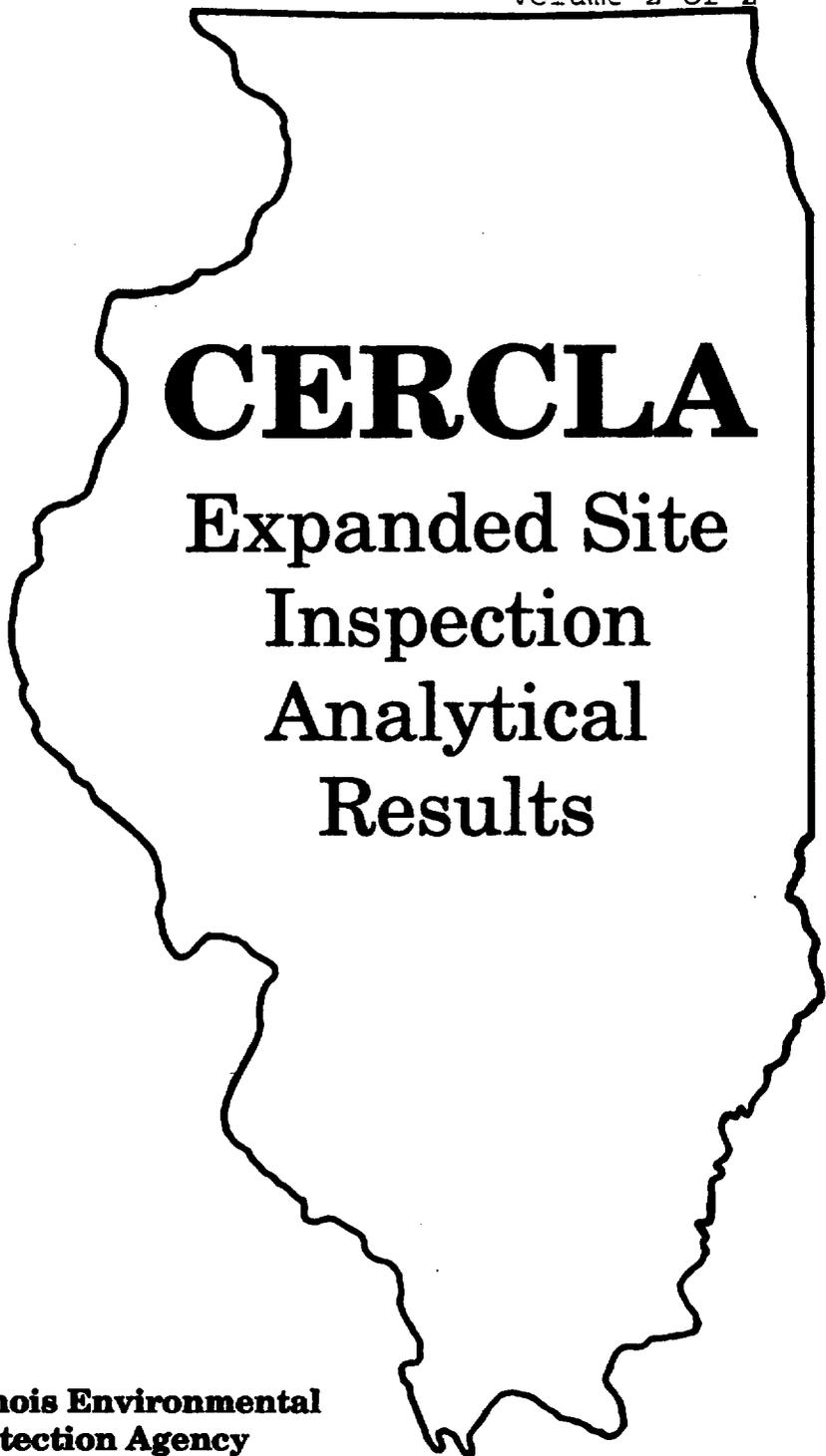


L1358070001 Montgomery Co.
Eagle Zinc Company
ILD 980606941
SF/HRS
Volume 2 of 2

LZ
1/17/96
156083



CERCLA

Expanded Site Inspection Analytical Results



**Illinois Environmental
Protection Agency**
P. O. Box 19276
Springfield, IL 62794-9276

Confidential Material May be Enclosed

APPENDIX J
ANALYTICAL RESULTS OF OCTOBER 26-27, 1993
EXPANDED SITE INSPECTION SAMPLES
ANALYZED FOR TARGET COMPOUND LIST PARAMETERS

DATA QUALIFIERS

QUALIFIER	DEFINITION ORGANICS	DEFINITION INORGANICS
U	Compound was tested for but not detected. The sample quantitation limit must be corrected for dilution and for percent moisture. For soil samples subjected to GPC clean-up procedures, the CRQL is also multiplied by two, to account for the fact that only half of the extract is recovered.	Analyte was analyzed for but not detected.
J	Estimated value. Used when estimating a concentration for tentatively identified compounds (TICS) where a 1:1 response is assumed or when the mass spectral data indicate the presence of a compound that meets the identification criteria and the result is less than the sample quantitation limit but greater than zero. Used in data validation when the quality control data indicate that a value may not be accurate.	Estimated value. Used in data validation when the quality control data indicate that a value may not be accurate.
C	This flag applies to pesticide results where the identification is confirmed by GC/MS.	Method qualifier indicates analysis by the Manual Spectrophotometric method.
B	Analyte was found in the associated blank as well as in the sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.	The reported value is less than the CRDL but greater than the instrument detection limit (IDL).
D	Identifies all compounds identified in an analysis at a secondary dilution factor. If a sample or extract is re-analyzed at a higher dilution factor as in the "E" flag, the "DL" suffix is appended to the sample number on the Form I for the diluted sample, and all concentration values are flagged with the "D" flag.	Not used.
E	Identifies compounds whose concentrations exceed the calibration range for that specific analysis. All extracts containing compounds exceeding the calibration range must be diluted and analyzed again. If the dilution of the extract causes any compounds identified in the first analysis to be below the calibration range in the second analysis, then the results of both analyses must be reported on separate Forms I. The Form I for the diluted sample must have the "DL" suffix appended to the sample number.	The reported value is estimated because of the presence of interference.
A	This flag indicates that a TIC is a suspected aldol concentration product formed by the reaction of the solvents used to process the sample in the laboratory.	Method qualifier indicates analysis by Flame Atomic Absorption (AA).
M	Not used.	Duplicate injection (a QC parameter not met).

N	Not used	Spiked sample (a QC parameter not met).
S	Not used.	The reported value was determined by the Method of Standard Additions (MSA).
W	Not used.	Post digestion spike for Furnace AA analysis (a QC parameter) is out of control limits of 85% to 115% recovery, while sample absorbance is less than 50% of spike absorbance.
*	Not used.	Duplicate analysis (a QC parameter not within control limits).
+	Not used.	Correlation coefficient for MSA (a QC parameter) is less than 0.995.
P	Not used.	Method qualifier indicates analysis by ICP (Inductively Coupled Plasma) Spectroscopy.
CV	Not used.	Method qualifier indicates analysis by Cold Vapor AA.
AV	Not used.	Method qualifier indicates analysis by Automated Cold Vapor AA.
AS	Not used.	Method qualifier indicates analysis by Semi-Automated Cold Spectrophotometry.
T	Not used.	Method qualifier indicates Titrimetric analysis.
NR	The analyte was not required to be analyzed.	The analyte was not required to be analyzed.
R	Rejected data. The QC parameters indicate that the data is not usable for any purpose.	Rejected data. The QC parameters indicate that the data is not usable for any purpose.

SITE NAME: EAGLE ZINC COMPANY

ILD 980606941

TABLE 3-2
SEDIMENT SUMMARY

SAMPLING POINT	X201	X202	X203	X204	X205	X206	X207	X208
PARAMETER	Backgd. Sediment	Dup of X201 Sediment	Sediment	Sediment	Sediment	Sediment	Sediment	Sediment
VOLATILES UG/KG								
Methylene Chloride	--	--	--	--	--	160.0 J	--	--
Acetone	11.0 J	22.0	12.0 J	22.0 UJ	37.0 J	76.0 J	--	17.0 UJ
2-Butanone (MEK)	14.0 UJ	4.0 J	6.0 J	22.0 UJ	20.0 J	48.0 J	14.0 UJ	17.0 UJ
1,1,1-Trichloroethane	--	--	17.0 UJ	27.0 J	9.0 J	290.0 J	--	8.0 J
Carbon Tetrachloride	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Bromodichloromethane	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
1,2-Dichloropropane	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
cis-1,3-Dichloropropane	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Trichloroethene	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Dibromochloromethane	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
1,1,2-Trichloroethane	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Benzene	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Trans-1,3-Dichloropropane	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Bromoform	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
4-Methyl-2-Pentanone	--	--	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
2-Hexanone	14.0 UJ	14.0 UJ	17.0 UJ	22.0 UJ	14.0 UJ	36.0 UJ	14.0 UJ	17.0 UJ
Tetrachloroethene	--	--	--	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Toluene	--	--	--	22.0 UJ	14.0 UJ	36.0 J	--	17.0 UJ
1,1,2,2-Tetrachloroethane	--	--	--	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Chlorobenzene	--	--	--	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Ethylbenzene	--	--	--	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Styrene	--	--	--	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
Xylene(total)	--	--	--	22.0 UJ	14.0 UJ	36.0 UJ	--	17.0 UJ
SEMIVOLATILES UG/KG								
4-Chloroaniline	470.0 UJ	470.0 UJ	560.0 UJ	730.0 UJ	480.0 UJ	1200.0 UJ	440.0 UJ	560.0 UJ
2-Methylnaphthalene	--	--	--	--	100.0 J	--	--	--
3-Nitroaniline	1100.0 UJ	1100.0 UJ	1400.0 UJ	1800.0 UJ	1200.0 UJ	2800.0 UJ	1100.0 UJ	1400.0 UJ
4-Nitroaniline	1100.0 R	1100.0 R	1400.0 R	1800.0 R	1200.0 R	2800.0 R	1100.0 R	1400.0 R
Phenanthrene	--	--	260.0 J	190.0	--	--	--	--
Anthracene	--	--	--	320.0 J	--	--	--	--
Carbazole	--	--	--	290.0 J	--	--	--	--
Fluoranthene	--	--	520.0 J	170.0	--	--	130.0 J	--
Pyrene	--	--	520.0 J	160.0	--	--	140.0 J	--
3,3'-Dichlorobenzidine	470.0 UJ	470.0 UJ	560.0 UJ	730.0 UJ	480.0 UJ	1200.0 UJ	440.0 UJ	560.0 UJ
Benzo(a)anthracene	--	--	230.0 J	650.0	--	--	100.0 J	--
Chrysene	--	--	310.0 J	670.0 J	--	--	120.0 J	--
Di(2-Ethylhexyl)phthalate	--	--	660.0	--	--	--	--	--
Benzo(b)fluoranthene	--	--	480.0 J	--	--	--	140.0 J	--
Benzo(k)fluoranthene	--	--	--	1200.0	--	--	--	--
Benzo(a)pyrene	--	--	230.0 J	810.0	--	--	--	--

SITE NAME: EAGLE ZINC COMPANY		TABLE 3-2 SEDIMENT SUMMARY							
ILD 980606941		X201	X202	X203	X204	X205	X206	X207	X208
SAMPLING POINT		Backgd.	Dup of X201						
PARAMETER		Sediment	Sediment	Sediment	Sediment	Sediment	Sediment	Sediment	Sediment
PESTICIDES UG/KG									
alpha-BHC		--	--	--	--	--	1.5 J	--	--
beta-BHC		--	--	--	--	--	1.0 JP	--	--
gamma-BHC (Lindane)		--	--	--	--	--	1.1 JP	--	--
Aldrin		--	--	4.4 P	--	--	--	--	--
Heptachlor epoxide		--	0.2 JP	--	1.9 JP	--	4.7 J	--	--
Dieldrin		2.3 J	2.6 J	16.0 P	12.0 P	--	10.0 J	--	1.3 JP
4,4'-DDE		--	0.4 JP	--	--	--	0.7 JP	--	--
Endrin		0.3 JP	0.9 J	18.0 P	12.0	2.4 J	--	--	2.8 JP
Endosulfan II		--	--	--	--	--	--	--	3.6 JP
4,4'-DDD		0.4 JP	0.9 JP	7.5 P	6.0 JP	--	1.8 JP	--	5.1 J
4,4'-DDT		3.7 J	0.4 J	11.0 P	15.0 P	--	4.8 J	--	--
Methoxychlor (Mariate)		--	--	--	--	--	13.0 J	--	--
Endrin Ketone		--	0.5 J	--	--	1.6 J	--	--	--
alpha-Chlorodane		2.0 JP	3.1 P	16.0 P	7.0 P	--	1.7 JP	--	0.6 J
gamma-Chlorodane		2.0 J	2.5	15.0 P	7.4 P	--	3.0 J	--	0.7 JP
Toxaphene		--	110.0 JP	--	--	--	--	--	320.0 P
Aroclor-1254		--	--	250.0	120.0	--	--	--	24.0 JP
Aroclor-1260		17.0 J	9.3 J	110.0 P	100.0	--	--	--	--
INORGANICS MG/KG									
Aluminum		6890.0	6390.0	7370.0	14900.0	8360.0	16300.0	10700.0	9810.0
Antimony		9.0 J	10.4 J	10.3 J	17.4 J	9.3 J	62.7 J	10.7 J	10.8 J
Arsenic		4.5	4.3	6.4	10.9	2.9	19.4	6.0	6.0
Barium		79.5	70.4	99.9	97.4	89.6	383.0	167.0	92.5
Beryllium		0.4 B	0.4 B	0.5 B	0.6 B	0.5 B	1.5 B	0.7 B	0.6 B
Cadmium		0.7 B	--	8.6	7.4	1.8	523.0	11.1	19.6
Calcium		6360.0	5520.0	20300.0	12000.0	4680.0	8280.0	1510.0	3020.0
Chromium		9.9	9.9	12.1	13.2	11.0	28.6	14.6	13.7
Cobalt		6.1 B	4.9 B	6.0 B	8.1 B	4.5 B	353.0	10.8 B	4.7
Copper		11.9	11.2	37.9	41.9	9.0	1420.0	20.8	52.2
Iron		10100.0	9120.0	12400.0	14300.0	10900.0	82400.0	14900.0	14500.0
Lead		46.4	35.0	101.0	72.6	10.2	932.0	76.0	125.0
Magnesium		2760.0	2390.0	3330.0	2960.0	2620.0	4970.0	1500.0	1930.0
Manganese		501.0	384.0	722.0	451.0	85.9	3500.0	1470.0	461.0
Mercury		--	--	0.2	0.1 B	--	0.7	--	0.3
Nickel		9.2 B	8.7 B	11.5	14.7 B	12.6	583.0	11.9	12.7
Selenium		0.3 J	0.3 J	0.3 J	0.4 J	0.3 J	4.1	0.3 J	0.4 J
Silver		0.2	--	--	--	--	14.1	--	--
Sodium		73.3 B	79.8 B	132.0 B	150.0 B	84.7 B	470.0 B	82.0 B	110.0 B
Thallium		0.3 J	--	--	0.4 J	0.3 J	3.8 J	0.3 J	0.4 J
Vanadium		17.9	17.4	19.0	26.3	20.8	52.9	41.2	27.2
Zinc		326.0	291.0	2200.0	3040.0	5690.0	156000.0	2410.0	3280.0

SITE NAME: EAGLE ZINC CO.

ILD 980606941

TABLE 3-2
SOIL SUMMARY

SAMPLING POINT PARAMETER	X101 Backgd. Soil	X102 Dup of X101 Soil	X103 Soil	X104 Soil	X105 Soil	X106 Soil
INORGANICS MG\KG (ppm)						
Aluminum	12400.00	10000.00	14900.00	6880.00	7430.00	13000.00
Antimony	8.90 J	9.20 J	13.90 J	10.60 J	11.40 J	9.40 J
Arsenic	5.80	5.70	5.00	6.60	86.30	6.20
Barium	230.00	265.00	112.00	181.00	379.00	224.00
Beryllium	0.80 B	0.81 B	0.68 B	0.49 B	0.83 B	0.63 B
Cadmium	--	--	3.20	3.20	47.20	0.89 B
Calcium	10600.00	9880.00	2010.00	598.00 B	1930.00	11600.00
Chromium	16.20	14.40	15.90	10.30	22.60	15.10
Cobalt	4.10 B	6.50 B	12.00 B	13.70	20.10	11.10
Copper	20.00 J	19.70 J	201.00 J	30.60 J	911.00 J	24.70 J
Iron	14700.00	14400.00	13900.00	11500.00	104000.00	15400.00
Lead	148.00	236.00	260.00	61.00	5760.00	28.50
Magnesium	2370.00	2090.00	1970.00	1040.00 B	1630.00	2150.00
Manganese	434.00	686.00	915.00	1180.00	178.00	922.00
Mercury	0.17	0.18	--	--	--	--
Nickel	13.50	11.50	20.00	27.10	55.90	14.00
Potassium	1890.00	1600.00	1120.00 B	491.00 J	300.00 J	1060.00 J
Selenium	--	1.30 J	0.31 J	0.27 J	1.30	--
Silver	--	--	--	--	6.30	--
Sodium	106.00 B	87.90 B	47.80 B	47.50 B	39.60 B	37.40 B
Thallium	0.33 B	0.34 J	0.31 J	1.20 J	1.30 J	0.26 J
Vanadium	28.50	27.10	28.20	27.50	22.60	28.50
Zinc	136.00	138.00	5580.00	4770.00	31700.00	1490.00

SITE NAME: EAGLE ZINC CO.

ILD 980606941

TABLE 3-2
SOIL SUMMARY

SAMPLING POINT	X107	X108	X109	X110	X111	X112
PARAMETER	Soil	Soil	Soil	Soil	Soil	Soil
INORGANICS MG\KG (ppm)						
Aluminum	13000.00	11500.00	10200.00	15000.00	13500.00	9950.00
Antimony	10.50 J	13.00 J	9.30 J	7.90 J	9.00 J	10.20 J
Arsenic	8.70	13.40	4.60	13.60	8.50	6.20
Barium	124.00	267.00	130.00	150.00	193.00	233.00
Beryllium	0.72 B	1.00 B	0.60 B	0.78 B	0.94 B	0.85 B
Cadmium	3.50	11.30	0.71 B	2.00	1.60	2.80
Calcium	5360.00	5430.00	2580.00	3450.00	8380.00	2800.00
Chromium	16.10	23.40	13.40	20.70	20.20	14.80
Cobalt	5.60 B	14.80	6.90 B	8.50 B	7.80 B	11.30 B
Copper	36.40 J	104.00	15.30	22.50	33.80	15.90
Iron	14900.00	33900.00	12600.00	20700.00	19600.00	13900.00
Lead	105.00	388.00	47.00	87.60	70.80	70.10
Magnesium	2090.00	1630.00	1530.00	2500.00	1950.00	1760.00
Manganese	600.00	1670.00	660.00	563.00	491.00	2070.00
Mercury	0.16	0.16	0.11 B	--	0.11 B	0.11 B
Nickel	15.90	35.10	11.00	15.90	16.50	22.90
Potassium	1160.00 J	--	1650.00	1980.00	1920.00	1970.00
Selenium	--	0.84 J	0.31 J	0.49 J	0.42 J	0.39 J
Silver	--	--	--	--	--	--
Sodium	71.80 B	178.00 B	65.70 B	62.80 B	120.00 B	52.40 B
Thallium	0.35 J	1.40 J	0.28 J	--	0.25 J	0.28 J
Vanadium	27.30	37.70	24.70	38.70	34.20	28.20
Zinc	2480.00	2280.00	360.00	606.00	488.00	489.00

SITE NAME: EAGLE ZINC CO.

ILD 980606941

SOIL TABLE 3-2
SUMMARY

SAMPLING POINT	X113	X114	X115	X116	X117
PARAMETER	Soil	Soil	Soil	Soil	Soil
INORGANICS MG\KG (ppm)					
Aluminum	16600.00	9750.00	14800.00	12500.00	13800.00
Antimony	7.80 J	8.40 J	11.10 J	9.90 J	14.50 J
Arsenic	5.60	11.90	10.50	7.10	8.50
Barium	116.00	183.00	181.00	227.00	222.00
Beryllium	0.85 B	1.00	0.80 B	0.93 B	1.70
Cadmium	0.68 B	2.90	1.48	2.30	4.80
Calcium	5940.00	4230.00	4970.00	8430.00	19300.00
Chromium	21.70	15.90	19.40	18.90	17.30
Cobalt	10.60	5.80 B	7.00 B	9.80 B	10.60 B
Copper	22.50	28.30 J	27.80 J	25.50 J	57.20 J
Iron	20400.00	28600.00	19700.00	18900.00	21100.00
Lead	75.10	137.00	76.20	147.00	186.00
Magnesium	4870.00	1130.00	2030.00	2020.00	2140.00
Manganese	568.00	314.00	538.00	851.00	995.00
Mercury	--	--	0.42	0.24	0.14 B
Nickel	18.60	14.40	10.90	16.50	27.50
Potassium	2400.00	1040.00	1470.00	1750.00	1460.00 J
Selenium	0.27 J	0.76 J	0.52 J	0.53 J	0.35 J
Silver	--	--	1.20	--	--
Sodium	45.80	293.00 B	61.50 B	89.90 B	1020.00 B
Thallium	0.27 J	0.71 J	0.57 J	0.53 J	0.35 J
Vanadium	33.70	29.70	34.80	35.10	34.30
Zinc	451.00	1580.00	638.00	998.00	7420.00

SITE NAME: EAGLE ZINC CO.

ILD 980606941

SOIL

TABLE 3-2
SUMMARY

SAMPLING POINT	X118	X119	X120
PARAMETER	Soil	Soil	Soil
INORGANICS MG\KG (ppm)			
Aluminum	14100.00	9390.00	16300.00
Antimony	10.90 J	8.30 J	8.00 J
Arsenic	5.90	6.70	10.70
Barium	106.00	196.00	155.00
Beryllium	0.73 B	0.60 B	0.95
Cadmium	--	2.80	--
Calcium	1720.00	12100.00	2870.00
Chromium	18.50	13.70	20.40
Cobalt	11.10 B	14.90	7.40 B
Copper	15.90 J	17.50 J	17.20 J
Iron	18200.00	14100.00	22900.00
Lead	30.40	51.90	32.70
Magnesium	2120.00	1790.00	2870.00
Manganese	795.00	1520.00	889.00
Mercury	--	0.32	--
Nickel	12.80	14.80	16.90
Potassium	1210.00 J	1670.00	1490.00
Selenium	0.27 J	0.55 J	0.38 J
Silver	--	--	--
Sodium	--	--	27.70 B
Thallium	0.27 J	0.50 J	0.25 J
Vanadium	34.50 B	26.70	39.00
Zinc	354.00	1570.00	371.00

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X201

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344846

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1029BK04

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 30 Date Analyzed: 10/29/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	14	U
74-83-9	-----Bromomethane	14	U
75-01-4	-----Vinyl Chloride	14	U
75-00-3	-----Chloroethane	14	U
75-09-2	-----Methylene Chloride	14 8	U <i>am</i>
67-64-1	-----Acetone	11	J
75-15-0	-----Carbon Disulfide	14	U
75-35-4	-----1,1-Dichloroethene	14	U
75-34-3	-----1,1-Dichloroethane	14	U
540-59-0	-----1,2-Dichloroethene (total)	14	U
107-06-2	-----1,2-Dichloroethane	14	U
78-93-3	-----2-Butanone	14	U
71-55-6	-----1,1,1-Trichloroethane	14	U
56-23-5	-----Carbon Tetrachloride	14	U
75-27-4	-----Bromodichloromethane	14	U
78-87-5	-----1,2-Dichloropropane	14	U
10061-01-5	-----cis-1,3-Dichloropropene	14	U
79-01-6	-----Trichloroethene	14	U
124-48-1	-----Dibromochloromethane	14	U
79-00-5	-----1,1,2-Trichloroethane	14	U
71-43-2	-----Benzene	14	U
10061-02-6	-----trans-1,3-Dichloropropene	14	U
75-25-2	-----Bromoform	14	U
108-10-1	-----4-Methyl-2-Pentanone	14	U
591-78-6	-----2-Hexanone	14	U
127-18-4	-----Tetrachloroethene	14	U
79-34-5	-----1,1,2,2-Tetrachloroethane	14	U
108-88-3	-----Toluene	14	U
108-90-7	-----Chlorobenzene	14	U
100-41-4	-----Ethylbenzene	14	U
100-42-5	-----Styrene	14	U
1330-20-7	-----Xylene (total)	14	U

RECEIVED

DEC 30 1993

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X201

Lab Name: ILLINOIS EPA Contract: 1358070001
Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846
Matrix: (soil/water) SOIL Lab Sample ID: D344846
Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1029BK04
Level: (low/med) LOW Date Received: 10/27/93
% Moisture: not dec. 30 Date Analyzed: 10/29/93
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X201

Lab Name: ILLINOIS EPA Contract: 1358070001
 Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846
 Matrix: (soil/water) SOIL Lab Sample ID: D344846
 Sample wt/vol: 30.50 (g/mL) G Lab File ID: B1102E09
 Level: (low/med) LOW Date Received: 10/27/93
 % Moisture: 31 decanted: (Y/N) N Date Extracted: 10/28/93
 Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0
 GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	470	U
111-44-4	bis(2-Chloroethyl) Ether	470	U
95-57-8	2-Chlorophenol	470	U
541-73-1	1,3-Dichlorobenzene	470	U
106-46-7	1,4-Dichlorobenzene	470	U
95-50-1	1,2-Dichlorobenzene	470	U
95-48-7	2-Methylphenol	470	U
108-60-1	2,2'-oxybis(1-Chloropropane)	470	U
106-44-5	4-Methylphenol	470	U
621-64-7	N-Nitroso-Di-n-Propylamine	470	U
67-72-1	Hexachloroethane	470	U
98-95-3	Nitrobenzene	470	U
78-59-1	Isophorone	470	U
88-75-5	2-Nitrophenol	470	U
105-67-9	2,4-Dimethylphenol	470	U
111-91-1	bis(2-Chloroethoxy)Methane	470	U
120-83-2	2,4-Dichlorophenol	470	U
120-82-1	1,2,4-Trichlorobenzene	470	U
91-20-3	Naphthalene	470	U
106-47-8	4-Chloroaniline	470	U ^J
87-68-3	Hexachlorobutadiene	470	U
59-50-7	4-Chloro-3-Methylphenol	470	U
91-57-6	2-Methylnaphthalene	470	U
77-47-4	Hexachlorocyclopentadiene	470	U
88-06-2	2,4,6-Trichlorophenol	470	U
95-95-4	2,4,5-Trichlorophenol	1100	U
91-58-7	2-Chloronaphthalene	470	U
88-74-4	2-Nitroaniline	1100	U
131-11-3	Dimethylphthalate	470	U
208-96-8	Acenaphthylene	470	U
606-20-2	2,6-Dinitrotoluene	470	U
99-09-2	3-Nitroaniline	1100	U ^J
83-32-9	Acenaphthene	470	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X201

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344846

Sample wt/vol: 30.50 (g/mL) G Lab File ID: B1102E09

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 31 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION	UNITS
51-28-5	2,4-Dinitrophenol	1100	U
100-02-7	4-Nitrophenol	1100	U
132-64-9	Dibenzofuran	470	U
121-14-2	2,4-Dinitrotoluene	470	U
84-66-2	Diethylphthalate	470	U
7005-72-3	4-Chlorophenyl-phenylether	470	U
86-73-7	Fluorene	470	U
100-10-6	4-Nitroaniline	1100	UR
534-52-1	4,6-Dinitro-2-methylphenol	1100	U
86-30-6	N-Nitrosodiphenylamine (1)	470	U
101-55-3	4-Bromophenyl-phenylether	470	U
118-74-1	Hexachlorobenzene	470	U
87-86-5	Pentachlorophenol	1100	U
85-01-8	Phenanthrene	470	U
120-12-7	Anthracene	470	U
86-74-8	Carbazole	470	U
84-74-2	Di-n-Butylphthalate	800	BU
206-44-0	Fluoranthene	470	U
129-00-0	Pyrene	470	U
85-68-7	Butylbenzylphthalate	470	U
91-94-1	3,3'-Dichlorobenzidine	470	UJ
56-55-3	Benzo(a)Anthracene	470	U
218-01-9	Chrysene	470	U
117-81-7	bis(2-Ethylhexyl) Phthalate	470	U
117-84-0	Di-n-Octyl Phthalate	470	U
205-99-2	Benzo(b)Fluoranthene	470	U
207-08-9	Benzo(k)Fluoranthene	470	U
50-32-8	Benzo(a)Pyrene	470	U
193-39-5	Indeno(1,2,3-cd)Pyrene	470	U
53-70-3	Dibenz(a,h)Anthracene	470	U
191-24-2	Benzo(g,h,i)Perylene	470	U

am

am

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X201

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344846

Sample wt/vol: 30.50 (g/mL) G Lab File ID: B1102E09

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 31 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

Number TICs found: 27

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q	
1.	UNKNOWN	10.17	1600	BSU	am
2.	UNKNOWN	10.88	650	BSU	am
3.	UNKNOWN	13.13	620	J	
4.	UNKNOWN ALIP. ACID	22.35	400	BSU	am
5.	UNKNOWN	25.50	130	J	
6.	UNKNOWN ALIP. ACID	26.83	320	J	
7.	UNKNOWN	28.43	260	J	
8.	UNKNOWN	28.68	430	J	
9.	UNKNOWN	32.30	220	J	
10.	UNKNOWN	32.50	1100	J	
11.	UNKNOWN	32.77	120	J	
12.	UNKNOWN	32.83	280	J	
13.	UNKNOWN	33.78	120	J	
14.	UNKNOWN ALIP. HYDROCARBON	34.38	1900	J	
15.	UNKNOWN	34.40	440	J	
16.	UNKNOWN	34.68	570	J	
17.	UNKNOWN	35.47	360	J	
18.	UNKNOWN	36.03	960	J	
19.	UNKNOWN ALIP. HYDROCARBON	36.63	1100	J	
20.	UNKNOWN ALIP. HYDROCARBON	36.78	5300	J	
21.	UNKNOWN ALIP. HYDROCARBON	36.93	970	J	
22.	UNKNOWN	38.93	230	J	
23.	UNKNOWN	39.17	710	J	
24.	UNKNOWN ALIP. HYDROCARBON	40.20	4700	J	
25.	UNKNOWN	42.03	1200	J	
26.	UNKNOWN	45.72	700	J	
27.	UNKNOWN	47.48	560	J	

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X201

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344846

Sample wt/vol: 30.1 (g/mL) G Lab File ID: _____

% Moisture: 30 decanted: (Y/N) N Date Received: 10/27/93

Extraction: (SepF/Cont/Sonc) SONC Date Extracted: 10/28/93

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/04/93

Injection Volume: 1.00 (uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) Y pH: 7.0 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6	alpha-BHC	2.4	U
319-85-7	beta-BHC	2.4	U
319-86-8	delta-BHC	2.4	U
58-89-9	gamma-BHC (Lindane)	2.4	U
76-44-8	Heptachlor	2.4	U
309-00-2	Aldrin	2.4 0.49	BFP U Am
1024-57-3	Heptachlor epoxide	2.4	U
959-98-8	Endosulfan I	2.4	U
60-57-1	Dieldrin	2.3	J
72-55-9	4,4'-DDE	4.7	U
72-20-8	Endrin	0.32	JP
33213-65-9	Endosulfan II	4.7	U
50-29-3	4,4'-DDD	0.37	JP
1031-07-8	Endosulfan sulfate	4.7	U
50-29-3	4,4'-DDT	3.7	J
72-43-5	Methoxychlor	24	U
53494-70-5	Endrin ketone	4.7	U
7421-36-3	Endrin aldehyde	4.7	U
5103-71-9	alpha-Chlordane	2.0	JP
5103-74-2	gamma-Chlordane	2.0	J
8001-35-2	Toxaphene	240	U
12674-11-2	Aroclor-1016	47	U
11104-28-2	Aroclor-1221	95	U
11141-16-5	Aroclor-1232	47	U
53469-21-9	Aroclor-1242	47	U
12672-29-6	Aroclor-1248	47	U
11097-69-1	Aroclor-1254	47	U
11096-82-5	Aroclor-1260	17	J

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X202

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344847

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1029BK05

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 30 Date Analyzed: 10/29/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	14	U
74-83-9	-----Bromomethane	14	U
75-01-4	-----Vinyl Chloride	14	U
75-00-3	-----Chloroethane	14	U
75-09-2	-----Methylene Chloride	147	BU am
67-64-1	-----Acetone	22	
75-15-0	-----Carbon Disulfide	14	U
75-35-4	-----1,1-Dichloroethene	14	U
75-34-3	-----1,1-Dichloroethane	14	U
540-59-0	-----1,2-Dichloroethene (total)	14	U
107-06-2	-----1,2-Dichloroethane	14	U
78-93-3	-----2-Butanone	4	J
71-55-6	-----1,1,1-Trichloroethane	14	U
56-23-5	-----Carbon Tetrachloride	14	U
75-27-4	-----Bromodichloromethane	14	U
78-87-5	-----1,2-Dichloropropane	14	U
10061-01-5	-----cis-1,3-Dichloropropene	14	U
79-01-6	-----Trichloroethene	14	U
124-48-1	-----Dibromochloromethane	14	U
79-00-5	-----1,1,2-Trichloroethane	14	U
71-43-2	-----Benzene	14	U
10061-02-6	-----trans-1,3-Dichloropropene	14	U
75-25-2	-----Bromoform	14	U
108-10-1	-----4-Methyl-2-Pentanone	14	U
591-78-6	-----2-Hexanone	14	UJ
127-18-4	-----Tetrachloroethene	14	U
79-34-5	-----1,1,2,2-Tetrachloroethane	14	U
108-88-3	-----Toluene	14	U
108-90-7	-----Chlorobenzene	14	U
100-41-4	-----Ethylbenzene	14	U
100-42-5	-----Styrene	14	U
1330-20-7	-----Xylene (total)	14	U

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X202

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344847

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1029BK05

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 30 Date Analyzed: 10/29/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X202

Lab Name: ILLINOIS EPA Contract: 1358070001
 Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846
 Matrix: (soil/water) SOIL Lab Sample ID: D344847
 Sample wt/vol: 30.10 (g/mL) G Lab File ID: B1102E08
 Level: (low/med) LOW Date Received: 10/27/93
 % Moisture: 30 decanted: (Y/N) N Date Extracted: 10/28/93
 Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0
 GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	470	U
111-44-4	bis(2-Chloroethyl) Ether	470	U
95-57-8	2-Chlorophenol	470	U
541-73-1	1,3-Dichlorobenzene	470	U
106-46-7	1,4-Dichlorobenzene	470	U
95-50-1	1,2-Dichlorobenzene	470	U
95-48-7	2-Methylphenol	470	U
108-60-1	2,2'-oxybis(1-Chloropropane)	470	U
106-44-5	4-Methylphenol	470	U
621-64-7	N-Nitroso-Di-n-Propylamine	470	U
67-72-1	Hexachloroethane	470	U
98-95-3	Nitrobenzene	470	U
78-59-1	Isophorone	470	U
88-75-5	2-Nitrophenol	470	U
105-67-9	2,4-Dimethylphenol	470	U
111-91-1	bis(2-Chloroethoxy)Methane	470	U
120-83-2	2,4-Dichlorophenol	470	U
120-82-1	1,2,4-Trichlorobenzene	470	U
91-20-3	Naphthalene	470	U
106-47-8	4-Chloroaniline	470	UJ
87-68-3	Hexachlorobutadiene	470	U
59-50-7	4-Chloro-3-Methylphenol	470	U
91-57-6	2-Methylnaphthalene	470	U
77-47-4	Hexachlorocyclopentadiene	470	U
88-06-2	2,4,6-Trichlorophenol	470	U
95-95-4	2,4,5-Trichlorophenol	1100	U
91-58-7	2-Chloronaphthalene	470	U
88-74-4	2-Nitroaniline	1100	U
131-11-3	Dimethylphthalate	470	U
208-96-8	Acenaphthylene	470	U
606-20-2	2,6-Dinitrotoluene	470	U
99-09-2	3-Nitroaniline	1100	UJ
83-32-9	Acenaphthene	470	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X202

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344847

Sample wt/vol: 30.10 (g/mL) G Lab File ID: B1102E08

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 30 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION	Q
51-28-5	2,4-Dinitrophenol	1100	U
100-02-7	4-Nitrophenol	1100	U
132-64-9	Dibenzofuran	470	U
121-14-2	2,4-Dinitrotoluene	470	U
84-66-2	Diethylphthalate	470	U
7005-72-3	4-Chlorophenyl-phenylether	470	U
86-73-7	Fluorene	470	U
100-10-6	4-Nitroaniline	1100	UR
534-52-1	4,6-Dinitro-2-methylphenol	1100	U
86-30-6	N-Nitrosodiphenylamine (1)	470	U
101-55-3	4-Bromophenyl-phenylether	470	U
118-74-1	Hexachlorobenzene	470	U
87-86-5	Pentachlorophenol	1100	U
85-01-8	Phenanthrene	470	U
120-12-7	Anthracene	470	U
86-74-8	Carbazole	470	U
84-74-2	Di-n-Butylphthalate	870	BU
206-44-0	Fluoranthene	470	U
129-00-0	Pyrene	470	U
85-68-7	Butylbenzylphthalate	470	U
91-94-1	3,3'-Dichlorobenzidine	470	UJ
56-55-3	Benzo(a)Anthracene	470	U
218-01-9	Chrysene	470	U
117-81-7	bis(2-Ethylhexyl)Phthalate	470	U
117-84-0	Di-n-Octyl Phthalate	470	U
205-99-2	Benzo(b)Fluoranthene	470	U
207-08-9	Benzo(k)Fluoranthene	470	U
50-32-8	Benzo(a)Pyrene	470	U
193-39-5	Indeno(1,2,3-cd)Pyrene	470	U
53-70-3	Dibenz(a,h)Anthracene	470	U
191-24-2	Benzo(g,h,i)Perylene	470	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X202

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344847

Sample wt/vol: 30.10 (g/mL) G Lab File ID: B1102E08

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 30 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Number TICs found: 25

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	10.18	1600	BJU <i>am</i>
2.	UNKNOWN	10.90	630	BJU <i>am</i>
3.	UNKNOWN	13.15	630	J
4.	UNKNOWN ALIP. ACID	22.37	460	BJU <i>am</i>
5.	UNKNOWN	25.52	180	J
6.	UNKNOWN ALIP. ACID	26.85	240	J
7.	UNKNOWN	28.70	190	J
8.	UNKNOWN	31.53	210	J
9.	UNKNOWN	32.32	200	J
10.	UNKNOWN	32.50	1200	J
11.	UNKNOWN	33.80	130	J
12.	UNKNOWN ALIP. HYDROCARBON	34.40	1800	J
13.	UNKNOWN	34.42	510	J
14.	UNKNOWN	34.47	400	J
15.	UNKNOWN	34.70	370	J
16.	UNKNOWN	35.48	300	J
17.	UNKNOWN	36.03	890	J
18.	UNKNOWN ALIP. HYDROCARBON	36.65	720	J
19.	UNKNOWN ALIP. HYDROCARBON	36.80	4800	J
20.	UNKNOWN	36.95	530	J
21.	UNKNOWN	39.17	690	J
22.	UNKNOWN ALIP. HYDROCARBON	39.97	820	J
23.	UNKNOWN ALIP. HYDROCARBON	40.20	4600	J
24.	UNKNOWN	42.05	900	J
25.	UNKNOWN	47.48	920	J

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X202

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344847

Sample wt/vol: 31.0 (g/mL) G Lab File ID: _____

% Moisture: 30 decanted: (Y/N) N Date Received: 10/27/93

Extraction: (SepF/Cont/Sonc) SONC Date Extracted: 10/28/93

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/04/93

Injection Volume: 1.00 (uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) Y pH: 7.0 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND Q

319-84-6-----alpha-BHC	2.4	U
319-85-7-----beta-BHC	2.4	U
319-86-8-----delta-BHC	2.4	U
58-89-9-----gamma-BHC (Lindane)	2.4	U
76-44-8-----Heptachlor	2.4	U
309-00-2-----Aldrin	2.4	0.36 BJPU <i>am</i>
1024-57-3-----Heptachlor epoxide	0.17	JP
959-98-8-----Endosulfan I	2.4	U
60-57-1-----Dieldrin	2.6	J
72-55-9-----4,4'-DDE	0.44	JP
72-20-8-----Endrin	0.90	J
33213-65-9-----Endosulfan II	4.6	U
50-29-3-----4,4'-DDD	0.92	JP
1031-07-8-----Endosulfan sulfate	4.6	U
50-29-3-----4,4'-DDT	0.44	J
72-43-5-----Methoxychlor	24	U
53494-70-5-----Endrin ketone	0.51	J
7421-36-3-----Endrin aldehyde	4.6	U
5103-71-9-----alpha-Chlordane	3.1	P
5103-74-2-----gamma-Chlordane	2.5	
8001-35-2-----Toxaphene	110	JP
12674-11-2-----Aroclor-1016	46	U
11104-28-2-----Aroclor-1221	93	U
11141-16-5-----Aroclor-1232	46	U
53469-21-9-----Aroclor-1242	46	U
12672-29-6-----Aroclor-1248	46	U
11097-69-1-----Aroclor-1254	46	U
11096-82-5-----Aroclor-1260	9.3	J

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X203

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344848

Sample wt/vol: 30.60 (g/mL) G Lab File ID: B1102E11

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 42 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	560	U
111-44-4	bis(2-Chloroethyl) Ether	560	U
95-57-8	2-Chlorophenol	560	U
541-73-1	1,3-Dichlorobenzene	560	U
106-46-7	1,4-Dichlorobenzene	560	U
95-50-1	1,2-Dichlorobenzene	560	U
95-48-7	2-Methylphenol	560	U
108-60-1	2,2'-oxybis(1-Chloropropane)	560	U
106-44-5	4-Methylphenol	560	U
621-64-7	N-Nitroso-Di-n-Propylamine	560	U
67-72-1	Hexachloroethane	560	U
98-95-3	Nitrobenzene	560	U
78-59-1	Isophorone	560	U
88-75-5	2-Nitrophenol	560	U
105-67-9	2,4-Dimethylphenol	560	U
111-91-1	bis(2-Chloroethoxy)Methane	560	U
120-83-2	2,4-Dichlorophenol	560	U
120-82-1	1,2,4-Trichlorobenzene	560	U
91-20-3	Naphthalene	560	U
106-47-8	4-Chloroaniline	560	UJ
87-68-3	Hexachlorobutadiene	560	U
59-50-7	4-Chloro-3-Methylphenol	560	U
91-57-6	2-Methylnaphthalene	560	U
77-47-4	Hexachlorocyclopentadiene	560	U
88-06-2	2,4,6-Trichlorophenol	560	U
95-95-4	2,4,5-Trichlorophenol	1400	U
91-58-7	2-Chloronaphthalene	560	U
88-74-4	2-Nitroaniline	1400	U
131-11-3	Dimethylphthalate	560	U
208-96-8	Acenaphthylene	560	U
606-20-2	2,6-Dinitrotoluene	560	U
99-09-2	3-Nitroaniline	1400	UJ
83-32-9	Acenaphthene	560	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X203

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344848

Sample wt/vol: 30.60 (g/mL) G Lab File ID: B1102E11

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 42 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
51-28-5	2,4-Dinitrophenol	1400	U
100-02-7	4-Nitrophenol	1400	U
132-64-9	Dibenzofuran	560	U
121-14-2	2,4-Dinitrotoluene	560	U
84-66-2	Diethylphthalate	560	U
7005-72-3	4-Chlorophenyl-phenylether	560	U
86-73-7	Fluorene	560	U
100-10-6	4-Nitroaniline	1400	U R
534-52-1	4,6-Dinitro-2-methylphenol	1400	U
86-30-6	N-Nitrosodiphenylamine (1)	560	U
101-55-3	4-Bromophenyl-phenylether	560	U
118-74-1	Hexachlorobenzene	560	U
87-86-5	Pentachlorophenol	1400	U
85-01-8	Phenanthrene	260	J
120-12-7	Anthracene	560	U
86-74-8	Carbazole	560	U
84-74-2	Di-n-Butylphthalate	1400	U U
206-44-0	Fluoranthene	520	J
129-00-0	Pyrene	520	J
85-68-7	Butylbenzylphthalate	560	U
91-94-1	3,3'-Dichlorobenzidine	560	UJ
56-55-3	Benzo(a)Anthracene	230	J
218-01-9	Chrysene	310	J
117-81-7	bis(2-Ethylhexyl)Phthalate	660	
117-84-0	Di-n-Octyl Phthalate	560	U
205-99-2	Benzo(b)Fluoranthene	480	J
207-08-9	Benzo(k)Fluoranthene	560	U
50-32-8	Benzo(a)Pyrene	230	J
193-39-5	Indeno(1,2,3-cd)Pyrene	560	U
53-70-3	Dibenz(a,h)Anthracene	560	U
191-24-2	Benzo(g,h,i)Perylene	560	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X203

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344848

Sample wt/vol: 30.60 (g/mL) G Lab File ID: B1102E11

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 42 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Number TICs found: 26

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	10.17	1700	BJU <i>am</i>
2.	UNKNOWN	31.27	140	J
3.	UNKNOWN	31.53	390	J
4.	UNKNOWN	31.88	170	J
5.	UNKNOWN ALIP. HYDROCARBON	32.50	1800	J
6.	UNKNOWN	32.85	110	J
7.	UNKNOWN	33.42	610	J
8.	UNKNOWN	33.80	310	J
9.	UNKNOWN	34.28	280	J
10.	UNKNOWN ALIP. HYDROCARBON	34.40	2800	J
11.	UNKNOWN	34.70	380	J
12.	UNKNOWN	35.52	600	J
13.	UNKNOWN	35.90	220	J
14.	UNKNOWN	36.07	1600	J
15.	UNKNOWN	36.58	700	J
16.	UNKNOWN ALIP. HYDROCARBON	36.65	1000	J
17.	UNKNOWN ALIP. HYDROCARBON	36.82	6900	J
18.	UNKNOWN	36.95	1500	J
19.	UNKNOWN	37.18	170	J
20.	UNKNOWN	37.50	150	J
21.	UNKNOWN	37.67	390	J
22.	UNKNOWN	37.70	180	J
23.	UNKNOWN	38.15	1100	J
24.	UNKNOWN	38.35	750	J
25.	UNKNOWN	39.20	1500	J
26.	UNKNOWN ALIP. HYDROCARBON	40.23	4700	J

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X203

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344848

Sample wt/vol: 30.1 (g/mL) G Lab File ID: _____

% Moisture: 42 decanted: (Y/N) N Date Received: 10/27/93

Extraction: (SepF/Cont/Sonc) SONC Date Extracted: 10/28/93

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/04/93

Injection Volume: 1.00 (uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) Y pH: 7.1 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6	alpha-BHC	2.9	U
319-85-7	beta-BHC	2.9	U
319-86-8	delta-BHC	2.9	U
58-89-9	gamma-BHC (Lindane)	2.9	U
76-44-8	Heptachlor	2.9	U
309-00-2	Aldrin	4.4	P
1024-57-3	Heptachlor epoxide	2.9	U
959-98-8	Endosulfan I	2.9	U
60-57-1	Dieldrin	16	P
72-55-9	4,4'-DDE	5.7	U
72-20-8	Endrin	18	P
33213-65-9	Endosulfan II	5.7	U
50-29-3	4,4'-DDD	7.5	P
1031-07-8	Endosulfan sulfate	5.7	U
50-29-3	4,4'-DDT	11	P
72-43-5	Methoxychlor	29	U
53494-70-5	Endrin ketone	5.7	U
7421-36-3	Endrin aldehyde	5.7	U
5103-71-9	alpha-Chlordane	16	P
5103-74-2	gamma-Chlordane	15	P
8001-35-2	Toxaphene	290	U
12674-11-2	Aroclor-1016	57	U
11104-28-2	Aroclor-1221	120	U
11141-16-5	Aroclor-1232	57	U
53469-21-9	Aroclor-1242	57	U
12672-29-6	Aroclor-1248	57	U
11097-69-1	Aroclor-1254	250	U
11096-82-5	Aroclor-1260	110	P

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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X203RE

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344848RE

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK04

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 42 Date Analyzed: 11/01/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	17	U
74-83-9-----	Bromomethane	17	U
75-01-4-----	Vinyl Chloride	17	U
75-00-3-----	Chloroethane	17	U
75-09-2-----	Methylene Chloride	17	U
67-64-1-----	Acetone	12	J
75-15-0-----	Carbon Disulfide	17	U
75-35-4-----	1,1-Dichloroethene	17	U
75-34-3-----	1,1-Dichloroethane	17	U
540-59-0-----	1,2-Dichloroethene (total)	17	U
107-06-2-----	1,2-Dichloroethane	17	U
78-93-3-----	2-Butanone	6	J
71-55-6-----	1,1,1-Trichloroethane	17	J
56-23-5-----	Carbon Tetrachloride	17	J
75-27-4-----	Bromodichloromethane	17	J
78-87-5-----	1,2-Dichloropropane	17	J
10061-01-5-----	cis-1,3-Dichloropropene	17	J
79-01-6-----	Trichloroethene	17	J
124-48-1-----	Dibromochloromethane	17	J
79-00-5-----	1,1,2-Trichloroethane	17	J
71-43-2-----	Benzene	17	J
10061-02-6-----	trans-1,3-Dichloropropene	17	J
75-25-2-----	Bromoform	17	J
108-10-1-----	4-Methyl-2-Pentanone	17	J
591-78-6-----	2-Hexanone	17	J
127-18-4-----	Tetrachloroethene	17	U
79-34-5-----	1,1,2,2-Tetrachloroethane	17	U
108-88-3-----	Toluene	17	U
108-90-7-----	Chlorobenzene	17	U
100-41-4-----	Ethylbenzene	17	U
100-42-5-----	Styrene	17	U
1330-20-7-----	Xylene (total)	17	U

17 ~~8~~ *BJU am*

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X203RE

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344848RE

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK04

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 42 Date Analyzed: 11/01/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X204

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344849

Sample wt/vol: 30.30 (g/mL) G Lab File ID: B1102E12

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 55 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.2

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	730	U
111-44-4	bis(2-Chloroethyl) Ether	730	U
95-57-8	2-Chlorophenol	730	U
541-73-1	1,3-Dichlorobenzene	730	U
106-46-7	1,4-Dichlorobenzene	730	U
95-50-1	1,2-Dichlorobenzene	730	U
95-48-7	2-Methylphenol	730	U
108-60-1	2,2'-oxybis(1-Chloropropane)	730	U
106-44-5	4-Methylphenol	730	U
621-64-7	N-Nitroso-Di-n-Propylamine	730	U
67-72-1	Hexachloroethane	730	U
98-95-3	Nitrobenzene	730	U
78-59-1	Isophorone	730	U
88-75-5	2-Nitrophenol	730	U
105-67-9	2,4-Dimethylphenol	730	U
111-91-1	bis(2-Chloroethoxy)Methane	730	U
120-83-2	2,4-Dichlorophenol	730	U
120-82-1	1,2,4-Trichlorobenzene	730	U
91-20-3	Naphthalene	730	U
106-47-8	4-Chloroaniline	730	UJ
87-68-3	Hexachlorobutadiene	730	U
59-50-7	4-Chloro-3-Methylphenol	730	U
91-57-6	2-Methylnaphthalene	730	U
77-47-4	Hexachlorocyclopentadiene	730	U
88-06-2	2,4,6-Trichlorophenol	730	U
95-95-4	2,4,5-Trichlorophenol	1800	U
91-58-7	2-Chloronaphthalene	730	U
88-74-4	2-Nitroaniline	1800	U
131-11-3	Dimethylphthalate	730	U
208-96-8	Acenaphthylene	730	U
606-20-2	2,6-Dinitrotoluene	730	U
99-09-2	3-Nitroaniline	1800	UJ
83-32-9	Acenaphthene	730	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X204

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344849

Sample wt/vol: 30.30 (g/mL) G Lab File ID: B1102E12

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 55 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.2

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND Q

51-28-5-----	2,4-Dinitrophenol	1800	U
100-02-7-----	4-Nitrophenol	1800	U
132-64-9-----	Dibenzofuran	730	U
121-14-2-----	2,4-Dinitrotoluene	730	U
84-66-2-----	Diethylphthalate	730	U
7005-72-3-----	4-Chlorophenyl-phenylether	730	U
86-73-7-----	Fluorene	730	U
100-10-6-----	4-Nitroaniline	1800	UR
534-52-1-----	4,6-Dinitro-2-methylphenol	1800	U
86-30-6-----	N-Nitrosodiphenylamine (1)	730	U
101-55-3-----	4-Bromophenyl-phenylether	730	U
118-74-1-----	Hexachlorobenzene	730	U
87-86-5-----	Pentachlorophenol	1800	U
85-01-8-----	Phenanthrene	1900	
120-12-7-----	Anthracene	320	J
86-74-8-----	Carbazole	290	J
84-74-2-----	Di-n-Butylphthalate	2000	JU
206-44-0-----	Fluoranthene	1700	
129-00-0-----	Pyrene	1600	
85-68-7-----	Butylbenzylphthalate	730	U
91-94-1-----	3,3'-Dichlorobenzidine	730	UJ
56-55-3-----	Benzo(a)Anthracene	850	
218-01-9-----	Chrysene	670	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate	730	U
117-84-0-----	Di-n-Octyl Phthalate	730	U
205-99-2-----	Benzo(b)Fluoranthene	730	U
207-08-9-----	Benzo(k)Fluoranthene	1200	
50-32-8-----	Benzo(a)Pyrene	810	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	730	U
53-70-3-----	Dibenz(a,h)Anthracene	730	U
191-24-2-----	Benzo(g,h,i)Perylene	730	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X204

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344849

Sample wt/vol: 30.30 (g/mL) G Lab File ID: B1102E12

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 55 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.2

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Number TICs found: 28

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	10.20	2800	<i>BU am</i>
2.	UNKNOWN ALIP. HYDROCARBON	18.98	270	J
3.	UNKNOWN ALIP. HYDROCARBON	20.85	740	J
4.	UNKNOWN	29.32	780	J
5.	UNKNOWN	30.42	670	J
6.	UNKNOWN	30.73	250	J
7.	UNKNOWN	31.00	310	J
8.	UNKNOWN ALIP. ACID ESTER	31.60	2500	J
9.	UNKNOWN	32.52	2600	J
10.	UNKNOWN	32.87	260	J
11.	UNKNOWN	33.45	600	J
12.	UNKNOWN	33.82	250	J
13.	UNKNOWN	34.18	310	J
14.	UNKNOWN ALIP. HYDROCARBON	34.42	3300	J
15.	UNKNOWN	34.47	980	J
16.	UNKNOWN	35.52	570	J
17.	UNKNOWN	35.65	220	J
18.	UNKNOWN	35.92	170	J
19.	UNKNOWN	36.08	1200	J
20.	UNKNOWN	36.62	1100	J
21.	UNKNOWN	36.63	1800	J
22.	UNKNOWN ALIP. HYDROCARBON	36.83	6900	J
23.	UNKNOWN	36.98	1300	J
24.	UNKNOWN	37.20	200	J
25.	UNKNOWN PNA	37.68	1400	J
26.	UNKNOWN	38.38	820	J
27.	UNKNOWN	39.22	660	J
28.	UNKNOWN ALIP. HYDROCARBON	40.25	4800	J

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X204

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344849

Sample wt/vol: 30.6 (g/mL) G Lab File ID: _____

% Moisture: 55 decanted: (Y/N) N Date Received: 10/27/93

Extraction: (SepF/Cont/Sonc) SONC Date Extracted: 10/28/93

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/04/93

Injection Volume: 1.00 (uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) Y pH: 7.2 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6	alpha-BHC	3.7	U
319-85-7	beta-BHC	3.7	U
319-86-8	delta-BHC	3.7	U
58-89-9	gamma-BHC (Lindane)	3.7	U
76-44-8	Heptachlor	3.7	U
309-00-2	Aldrin	3.7	U
1024-57-3	Heptachlor epoxide	3.7 0.50	BJP U <i>am</i>
959-98-8	Endosulfan I	1.3	JP
60-57-1	Dieldrin	3.7	U
72-55-9	4,4'-DDE	12	P
72-20-8	Endrin	7.2	U
33213-65-9	Endosulfan II	12	U
50-29-3	4,4'-DDD	7.2	U
1031-07-8	Endosulfan sulfate	6.0	JP
50-29-3	4,4'-DDT	7.2	U
72-43-5	Methoxychlor	15	P
53494-70-5	Endrin ketone	37	U
7421-36-3	Endrin aldehyde	7.2	U
5103-71-9	alpha-Chlordane	7.2	U
5103-74-2	gamma-Chlordane	7.0	P
8001-35-2	Toxaphene	7.4	P
12674-11-2	Aroclor-1016	370	U
11104-28-2	Aroclor-1221	72	U
11141-16-5	Aroclor-1232	150	U
53469-21-9	Aroclor-1242	72	U
12672-29-6	Aroclor-1248	72	U
11097-69-1	Aroclor-1254	72	U
11096-82-5	Aroclor-1260	120	U
		100	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X204RE

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344849RE

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK05

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 55 Date Analyzed: 11/01/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	22	U
74-83-9	-----Bromomethane	22	U
75-01-4	-----Vinyl Chloride	22	U
75-00-3	-----Chloroethane	22	U
75-09-2	-----Methylene Chloride	22, 8	U <i>am</i>
67-64-1	-----Acetone	22	UJ
75-15-0	-----Carbon Disulfide	22	U
75-35-4	-----1,1-Dichloroethene	22	U
75-34-3	-----1,1-Dichloroethane	22	U
540-59-0	-----1,2-Dichloroethene (total)	22	U
107-06-2	-----1,2-Dichloroethane	22	U
78-93-3	-----2-Butanone	22	UJ
71-55-6	-----1,1,1-Trichloroethane	27	J
56-23-5	-----Carbon Tetrachloride	22	UJ
75-27-4	-----Bromodichloromethane	22	UJ
78-87-5	-----1,2-Dichloropropane	22	UJ
10061-01-5	-----cis-1,3-Dichloropropene	22	UJ
79-01-6	-----Trichloroethene	22	UJ
124-48-1	-----Dibromochloromethane	22	UJ
79-00-5	-----1,1,2-Trichloroethane	22	UJ
71-43-2	-----Benzene	22	UJ
10061-02-6	-----trans-1,3-Dichloropropene	22	UJ
75-25-2	-----Bromoform	22	UJ
108-10-1	-----4-Methyl-2-Pentanone	22	UJ
591-78-6	-----2-Hexanone	22	UJ
127-18-4	-----Tetrachloroethene	22	UJ
79-34-5	-----1,1,2,2-Tetrachloroethane	22	UJ
108-88-3	-----Toluene	22	UJ
108-90-7	-----Chlorobenzene	22	UJ
100-41-4	-----Ethylbenzene	22	UJ
100-42-5	-----Styrene	22	UJ
1330-20-7	-----Xylene (total)	22	UJ

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X204RE

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344849RE

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK05

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 55 Date Analyzed: 11/01/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X205

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344850

Sample wt/vol: 30.20 (g/mL) G Lab File ID: B1102E07

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 31 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND		
108-95-2	Phenol	480	U
111-44-4	bis(2-Chloroethyl) Ether	480	U
95-57-8	2-Chlorophenol	480	U
541-73-1	1,3-Dichlorobenzene	480	U
106-46-7	1,4-Dichlorobenzene	480	U
95-50-1	1,2-Dichlorobenzene	480	U
95-48-7	2-Methylphenol	480	U
108-60-1	2,2'-oxybis(1-Chloropropane)	480	U
106-44-5	4-Methylphenol	480	U
621-64-7	N-Nitroso-Di-n-Propylamine	480	U
67-72-1	Hexachloroethane	480	U
98-95-3	Nitrobenzene	480	U
78-59-1	Isophorone	480	U
88-75-5	2-Nitrophenol	480	U
105-67-9	2,4-Dimethylphenol	480	U
111-91-1	bis(2-Chloroethoxy)Methane	480	U
120-83-2	2,4-Dichlorophenol	480	U
120-82-1	1,2,4-Trichlorobenzene	480	U
91-20-3	Naphthalene	480	U
106-47-8	4-Chloroaniline	480	UJ
87-68-3	Hexachlorobutadiene	480	U
59-50-7	4-Chloro-3-Methylphenol	480	U
91-57-6	2-Methylnaphthalene	100	J
77-47-4	Hexachlorocyclopentadiene	480	U
88-06-2	2,4,6-Trichlorophenol	480	U
95-95-4	2,4,5-Trichlorophenol	1200	U
91-58-7	2-Chloronaphthalene	480	U
88-74-4	2-Nitroaniline	1200	U
131-11-3	Dimethylphthalate	480	U
208-96-8	Acenaphthylene	480	U
606-20-2	2,6-Dinitrotoluene	480	U
99-09-2	3-Nitroaniline	1200	UJ
83-32-9	Acenaphthene	480	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X205

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344850

Sample wt/vol: 30.20 (g/mL) G Lab File ID: B1102E07

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 31 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
51-28-5-----	2,4-Dinitrophenol	1200	U
100-02-7-----	4-Nitrophenol	1200	U
132-64-9-----	Dibenzofuran	480	U
121-14-2-----	2,4-Dinitrotoluene	480	U
84-66-2-----	Diethylphthalate	480	U
7005-72-3-----	4-Chlorophenyl-phenylether	480	U
86-73-7-----	Fluorene	480	U
100-10-6-----	4-Nitroaniline	1200	UR
534-52-1-----	4,6-Dinitro-2-methylphenol	1200	U
86-30-6-----	N-Nitrosodiphenylamine (1)	480	U
101-55-3-----	4-Bromophenyl-phenylether	480	U
118-74-1-----	Hexachlorobenzene	480	U
87-86-5-----	Pentachlorophenol	1200	U
85-01-8-----	Phenanthrene	480	U
120-12-7-----	Anthracene	480	U
86-74-8-----	Carbazole	480	U
84-74-2-----	Di-n-Butylphthalate	1200	BU
206-44-0-----	Fluoranthene	480	U
129-00-0-----	Pyrene	480	U
85-68-7-----	Butylbenzylphthalate	480	U
91-94-1-----	3,3'-Dichlorobenzidine	480	UJ
56-55-3-----	Benzo(a)Anthracene	480	U
218-01-9-----	Chrysene	480	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	480	U
117-84-0-----	Di-n-Octyl Phthalate	480	U
205-99-2-----	Benzo(b) Fluoranthene	480	U
207-08-9-----	Benzo(k) Fluoranthene	480	U
50-32-8-----	Benzo(a) Pyrene	480	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	480	U
53-70-3-----	Dibenz(a,h) Anthracene	480	U
191-24-2-----	Benzo(g,h,i) Perylene	480	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X205

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344850

Sample wt/vol: 30.20 (g/mL) G Lab File ID: B1102E07

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 31 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Number TICs found: 16

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q	
1.	UNKNOWN	10.18	1600	BJU	<i>an</i>
2.	UNKNOWN	10.90	540	BJU	<i>an</i>
3.	UNKNOWN	13.15	330	J	
4.	UNKNOWN	20.30	150	J	
5.	UNKNOWN ALIP. ACID	22.37	470	BJU	<i>an</i>
6.	UNKNOWN	23.63	200	J	
7.	UNKNOWN ALIP. HYDROCARBON	23.70	610	J	
8.	UNKNOWN	25.50	190	J	
9.	UNKNOWN ALIP. HYDROCARBON	26.13	180	J	
10.	UNKNOWN	32.50	660	J	
11.	UNKNOWN ALIP. HYDROCARBON	34.38	530	J	
12.	UNKNOWN	36.03	450	J	
13.	UNKNOWN ALIP. HYDROCARBON	36.78	1500	J	
14.	UNKNOWN	36.95	650	J	
15.	UNKNOWN ALIP. HYDROCARBON	40.18	1100	J	
16.	UNKNOWN	45.33	1200	J	

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X205

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344850

Sample wt/vol: 30.0 (g/mL) G Lab File ID: _____

% Moisture: 31 decanted: (Y/N) N Date Received: 10/27/93

Extraction: (SepF/Cont/Sonc) SONC Date Extracted: 10/28/93

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/04/93

Injection Volume: 1.00 (uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) Y pH: 7.0 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

319-84-6-----alpha-BHC	2.5	U
319-85-7-----beta-BHC	2.5	U
319-86-8-----delta-BHC	2.5	U
58-89-9-----gamma-BHC (Lindane)	2.5	U
76-44-8-----Heptachlor	2.5	U
309-00-2-----Aldrin	2.5	U
1024-57-3-----Heptachlor epoxide	2.5	U
959-98-8-----Endosulfan I	2.5	U
60-57-1-----Dieldrin	4.8	U
72-55-9-----4,4'-DDE	4.8	U
72-20-8-----Endrin	2.4	J
33213-65-9-----Endosulfan II	4.8	U
50-29-3-----4,4'-DDD	4.8	U
1031-07-8-----Endosulfan sulfate	4.8	U
50-29-3-----4,4'-DDT	4.8	U
72-43-5-----Methoxychlor	25	U
53494-70-5-----Endrin ketone	1.6	J
7421-36-3-----Endrin aldehyde	4.8	U
5103-71-9-----alpha-Chlordane	2.5	U
5103-74-2-----gamma-Chlordane	2.5	U
8001-35-2-----Toxaphene	250	U
12674-11-2-----Aroclor-1016	48	U
11104-28-2-----Aroclor-1221	97	U
11141-16-5-----Aroclor-1232	48	U
53469-21-9-----Aroclor-1242	48	U
12672-29-6-----Aroclor-1248	48	U
11097-69-1-----Aroclor-1254	48	U
11096-82-5-----Aroclor-1260	48	U

2.5 ~~0.28~~ BJPU am

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X205RE

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344850RE

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK06

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 31 Date Analyzed: 11/01/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	14	U
74-83-9	-----Bromomethane	14	U
75-01-4	-----Vinyl Chloride	14	U
75-00-3	-----Chloroethane	14	U
75-09-2	-----Methylene Chloride	24	U
67-64-1	-----Acetone	37	J
75-15-0	-----Carbon Disulfide	14	U
75-35-4	-----1,1-Dichloroethene	14	U
75-34-3	-----1,1-Dichloroethane	14	U
540-59-0	-----1,2-Dichloroethene (total)	14	U
107-06-2	-----1,2-Dichloroethane	14	U
78-93-3	-----2-Butanone	20	J
71-55-6	-----1,1,1-Trichloroethane	9	J
56-23-5	-----Carbon Tetrachloride	14	UJ
75-27-4	-----Bromodichloromethane	14	UJ
78-87-5	-----1,2-Dichloropropane	14	UJ
10061-01-5	-----cis-1,3-Dichloropropene	14	UJ
79-01-6	-----Trichloroethene	14	UJ
124-48-1	-----Dibromochloromethane	14	UJ
79-00-5	-----1,1,2-Trichloroethane	14	UJ
71-43-2	-----Benzene	14	UJ
10061-02-6	-----trans-1,3-Dichloropropene	14	UJ
75-25-2	-----Bromoform	14	UJ
108-10-1	-----4-Methyl-2-Pentanone	14	UJ
591-78-6	-----2-Hexanone	14	UJ
127-18-4	-----Tetrachloroethene	14	UJ
79-34-5	-----1,1,2,2-Tetrachloroethane	14	UJ
108-88-3	-----Toluene	14	UJ
108-90-7	-----Chlorobenzene	14	UJ
100-41-4	-----Ethylbenzene	14	UJ
100-42-5	-----Styrene	14	UJ
1330-20-7	-----Xylene (total)	14	UJ

am

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X205RE

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344850RE

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK06

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 31 Date Analyzed: 11/01/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____(uL)

Number TICs found: 1 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 629-14-1	ETHANE, 1,2-DIETHOXY-	6.45	17	JN

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X206

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344851

Sample wt/vol: 30.60 (g/mL) G Lab File ID: B1102E05

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 72 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/02/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	1200	U
111-44-4	bis(2-Chloroethyl) Ether	1200	U
95-57-8	2-Chlorophenol	1200	U
541-73-1	1,3-Dichlorobenzene	1200	U
106-46-7	1,4-Dichlorobenzene	1200	U
95-50-1	1,2-Dichlorobenzene	1200	U
95-48-7	2-Methylphenol	1200	U
108-60-1	2,2'-oxybis(1-Chloropropane)	1200	U
106-44-5	4-Methylphenol	1200	U
621-64-7	N-Nitroso-Di-n-Propylamine	1200	U
67-72-1	Hexachloroethane	1200	U
98-95-3	Nitrobenzene	1200	U
78-59-1	Isophorone	1200	U
88-75-5	2-Nitrophenol	1200	U
105-67-9	2,4-Dimethylphenol	1200	U
111-91-1	bis(2-Chloroethoxy)Methane	1200	U
120-83-2	2,4-Dichlorophenol	1200	U
120-82-1	1,2,4-Trichlorobenzene	1200	U
91-20-3	Naphthalene	1200	U
106-47-8	4-Chloroaniline	1200	UJ
87-68-3	Hexachlorobutadiene	1200	U
59-50-7	4-Chloro-3-Methylphenol	1200	U
91-57-6	2-Methylnaphthalene	1200	U
77-47-4	Hexachlorocyclopentadiene	1200	U
88-06-2	2,4,6-Trichlorophenol	1200	U
95-95-4	2,4,5-Trichlorophenol	2800	U
91-58-7	2-Chloronaphthalene	1200	U
88-74-4	2-Nitroaniline	2800	U
131-11-3	Dimethylphthalate	1200	U
208-96-8	Acenaphthylene	1200	U
606-20-2	2,6-Dinitrotoluene	1200	U
99-09-2	3-Nitroaniline	2800	UJ
83-32-9	Acenaphthene	1200	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X206

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344851

Sample wt/vol: 30.60 (g/mL) G Lab File ID: B1102E05

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 72 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/02/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
51-28-5-----	2,4-Dinitrophenol_____	2800	U
100-02-7-----	4-Nitrophenol_____	2800	U
132-64-9-----	Dibenzofuran_____	1200	U
121-14-2-----	2,4-Dinitrotoluene_____	1200	U
84-66-2-----	Diethylphthalate_____	1200	U
7005-72-3-----	4-Chlorophenyl-phenylether_____	1200	U
86-73-7-----	Fluorene_____	1200	U
100-10-6-----	4-Nitroaniline_____	2800	U R
534-52-1-----	4,6-Dinitro-2-methylphenol_____	2800	U
86-30-6-----	N-Nitrosodiphenylamine (1)_____	1200	U
101-55-3-----	4-Bromophenyl-phenylether_____	1200	U
118-74-1-----	Hexachlorobenzene_____	1200	U
87-86-5-----	Pentachlorophenol_____	2800	U
85-01-8-----	Phenanthrene_____	1200	U
120-12-7-----	Anthracene_____	1200	U
86-74-8-----	Carbazole_____	1200	U
84-74-2-----	Di-n-Butylphthalate_____	3200	U U
206-44-0-----	Fluoranthene_____	1200	U
129-00-0-----	Pyrene_____	1200	U
85-68-7-----	Butylbenzylphthalate_____	1200	U
91-94-1-----	3,3'-Dichlorobenzidine_____	1200	UJ
56-55-3-----	Benzo(a)Anthracene_____	1200	U
218-01-9-----	Chrysene_____	1200	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate_____	1200	U
117-84-0-----	Di-n-Octyl Phthalate_____	1200	U
205-99-2-----	Benzo(b)Fluoranthene_____	1200	U
207-08-9-----	Benzo(k)Fluoranthene_____	1200	U
50-32-8-----	Benzo(a)Pyrene_____	1200	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene_____	1200	U
53-70-3-----	Dibenz(a,h)Anthracene_____	1200	U
191-24-2-----	Benzo(g,h,i)Perylene_____	1200	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X206

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344851

Sample wt/vol: 30.60 (g/mL) G Lab File ID: B1102E05

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 72 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/02/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.1

Number TICs found: 20

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q	
1.	UNKNOWN	10.20	3800	BT U	am
2.	UNKNOWN	10.92	1100	BT U	am
3.	UNKNOWN ALIP. HYDROCARBON	20.85	530	J	
4.	UNKNOWN ALIP. ACID	22.38	1000	BT U	am
5.	UNKNOWN ALIP. ACID	26.87	1100	J	
6.	UNKNOWN	29.88	950	J	
7.	UNKNOWN	30.57	1800	J	
8.	UNKNOWN	31.55	1300	J	
9.	UNKNOWN	31.92	640	J	
10.	UNKNOWN	32.53	4400	J	
11.	UNKNOWN	32.68	360	J	
12.	UNKNOWN ALIP. HYDROCARBON	33.45	950	J	
13.	UNKNOWN	33.82	250	J	
14.	UNKNOWN	34.42	3100	J	
15.	UNKNOWN	34.50	1200	J	
16.	UNKNOWN	35.52	610	J	
17.	UNKNOWN	36.07	1900	J	
18.	UNKNOWN	36.68	390	J	
19.	UNKNOWN ALIP. HYDROCARBON	36.82	3600	J	
20.	UNKNOWN	47.48	4200	J	

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X206

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344851

Sample wt/vol: 30.8 (g/mL) G Lab File ID: _____

% Moisture: 72 decanted: (Y/N) N Date Received: 10/27/93

Extraction: (SepF/Cont/Sonc) SONC Date Extracted: 10/28/93

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/04/93

Injection Volume: 1.00 (uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) Y pH: 7.1 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6	alpha-BHC	1.5	J
319-85-7	beta-BHC	1.0	JP
319-86-8	delta-BHC	5.9	U
58-89-9	gamma-BHC (Lindane)	1.1	JP
76-44-8	Heptachlor	5.9	U
309-00-2	Aldrin	5.9 0.92	BJP U <i>am</i>
1024-57-3	Heptachlor epoxide	4.7	J
959-98-8	Endosulfan I	5.9	U
60-57-1	Dieldrin	10	J
72-55-9	4,4'-DDE	0.72	JP
72-20-8	Endrin	11	U
33213-65-9	Endosulfan II	11	U
50-29-3	4,4'-DDD	1.8	JP
1031-07-8	Endosulfan sulfate	11	U
50-29-3	4,4'-DDT	4.8	J
72-43-5	Methoxychlor	13	J
53494-70-5	Endrin ketone	11	U
7421-36-3	Endrin aldehyde	11	U
5103-71-9	alpha-Chlordane	1.7	JP
5103-74-2	gamma-Chlordane	3.0	J
8001-35-2	Toxaphene	590	U
12674-11-2	Aroclor-1016	110	U
11104-28-2	Aroclor-1221	230	U
11141-16-5	Aroclor-1232	110	U
53469-21-9	Aroclor-1242	110	U
12672-29-6	Aroclor-1248	110	U
11097-69-1	Aroclor-1254	110	U
11096-82-5	Aroclor-1260	110	U

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X206RE

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344851RE

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK07

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 72 Date Analyzed: 11/01/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	36	U
74-83-9	-----Bromomethane	36	U
75-01-4	-----Vinyl Chloride	36	U
75-00-3	-----Chloroethane	36	U
75-09-2	-----Methylene Chloride	160	J
67-64-1	-----Acetone	76	J
75-15-0	-----Carbon Disulfide	36	U
75-35-4	-----1,1-Dichloroethene	36	U
75-34-3	-----1,1-Dichloroethane	36	U
540-59-0	-----1,2-Dichloroethene (total)	36	U
107-06-2	-----1,2-Dichloroethane	36	U
78-93-3	-----2-Butanone	48	J
71-55-6	-----1,1,1-Trichloroethane	290	J
56-23-5	-----Carbon Tetrachloride	36	J
75-27-4	-----Bromodichloromethane	36	J
78-87-5	-----1,2-Dichloropropane	36	J
10061-01-5	-----cis-1,3-Dichloropropene	36	J
79-01-6	-----Trichloroethene	36	J
124-48-1	-----Dibromochloromethane	36	J
79-00-5	-----1,1,2-Trichloroethane	36	J
71-43-2	-----Benzene	36	J
10061-02-6	-----trans-1,3-Dichloropropene	36	J
75-25-2	-----Bromoform	36	J
108-10-1	-----4-Methyl-2-Pentanone	36	J
591-78-6	-----2-Hexanone	36	J
127-18-4	-----Tetrachloroethene	36	J
79-34-5	-----1,1,2,2-Tetrachloroethane	36	J
108-88-3	-----Toluene	36	J
108-90-7	-----Chlorobenzene	36	J
100-41-4	-----Ethylbenzene	36	J
100-42-5	-----Styrene	36	J
1330-20-7	-----Xylene (total)	36	J

am

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X206RE

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344851RE

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK07

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 72 Date Analyzed: 11/01/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Number TICs found: 2

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	6.48	74	J
2. 624-92-0	DISULFIDE, DIMETHYL	16.00	70	JN

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X207

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344852

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1029BK10

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 26 Date Analyzed: 10/29/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	14	U
74-83-9	-----Bromomethane	14	U
75-01-4	-----Vinyl Chloride	14	U
75-00-3	-----Chloroethane	14	U
75-09-2	-----Methylene Chloride	14	U
67-64-1	-----Acetone	14	U
75-15-0	-----Carbon Disulfide	14	U
75-35-4	-----1,1-Dichloroethene	14	U
75-34-3	-----1,1-Dichloroethane	14	U
540-59-0	-----1,2-Dichloroethene (total)	14	U
107-06-2	-----1,2-Dichloroethane	14	U
78-93-3	-----2-Butanone	14	UJ
71-55-6	-----1,1,1-Trichloroethane	14	U
56-23-5	-----Carbon Tetrachloride	14	U
75-27-4	-----Bromodichloromethane	14	U
78-87-5	-----1,2-Dichloropropane	14	U
10061-01-5	-----cis-1,3-Dichloropropene	14	U
79-01-6	-----Trichloroethene	14	U
124-48-1	-----Dibromochloromethane	14	U
79-00-5	-----1,1,2-Trichloroethane	14	U
71-43-2	-----Benzene	14	U
10061-02-6	-----trans-1,3-Dichloropropene	14	U
75-25-2	-----Bromoform	14	U
108-10-1	-----4-Methyl-2-Pentanone	14	U
591-78-6	-----2-Hexanone	14	UJ
127-18-4	-----Tetrachloroethene	14	U
79-34-5	-----1,1,2,2-Tetrachloroethane	14	U
108-88-3	-----Toluene	14	U
108-90-7	-----Chlorobenzene	14	U
100-41-4	-----Ethylbenzene	14	U
100-42-5	-----Styrene	14	U
1330-20-7	-----Xylene (total)	14	U

14 / ~~BJU~~ am

1E
 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X207

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344852

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1029BK10

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 26 Date Analyzed: 10/29/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X207

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344852

Sample wt/vol: 30.60 (g/mL) G Lab File ID: B1102E10

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 26 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND		
108-95-2	Phenol	440	U
111-44-4	bis(2-Chloroethyl) Ether	440	U
95-57-8	2-Chlorophenol	440	U
541-73-1	1,3-Dichlorobenzene	440	U
106-46-7	1,4-Dichlorobenzene	440	U
95-50-1	1,2-Dichlorobenzene	440	U
95-48-7	2-Methylphenol	440	U
108-60-1	2,2'-oxybis(1-Chloropropane)	440	U
106-44-5	4-Methylphenol	440	U
621-64-7	N-Nitroso-Di-n-Propylamine	440	U
67-72-1	Hexachloroethane	440	U
98-95-3	Nitrobenzene	440	U
78-59-1	Isophorone	440	U
88-75-5	2-Nitrophenol	440	U
105-67-9	2,4-Dimethylphenol	440	U
111-91-1	bis(2-Chloroethoxy)Methane	440	U
120-83-2	2,4-Dichlorophenol	440	U
120-82-1	1,2,4-Trichlorobenzene	440	U
91-20-3	Naphthalene	440	U
106-47-8	4-Chloroaniline	440	UJ
87-68-3	Hexachlorobutadiene	440	U
59-50-7	4-Chloro-3-Methylphenol	440	U
91-57-6	2-Methylnaphthalene	440	U
77-47-4	Hexachlorocyclopentadiene	440	U
88-06-2	2,4,6-Trichlorophenol	440	U
95-95-4	2,4,5-Trichlorophenol	1100	U
91-58-7	2-Chloronaphthalene	440	U
88-74-4	2-Nitroaniline	1100	U
131-11-3	Dimethylphthalate	440	U
208-96-8	Acenaphthylene	440	U
606-20-2	2,6-Dinitrotoluene	440	U
99-09-2	3-Nitroaniline	1100	UJ
83-32-9	Acenaphthene	440	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X207

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344852

Sample wt/vol: 30.60 (g/mL) G Lab File ID: B1102E10

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 26 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
51-28-5-----	2,4-Dinitrophenol	1100	U
100-02-7-----	4-Nitrophenol	1100	U
132-64-9-----	Dibenzofuran	440	U
121-14-2-----	2,4-Dinitrotoluene	440	U
84-66-2-----	Diethylphthalate	440	U
7005-72-3-----	4-Chlorophenyl-phenylether	440	U
86-73-7-----	Fluorene	440	U
100-10-6-----	4-Nitroaniline	1100	U R
534-52-1-----	4,6-Dinitro-2-methylphenol	1100	U
86-30-6-----	N-Nitrosodiphenylamine (1)	440	U
101-55-3-----	4-Bromophenyl-phenylether	440	U
118-74-1-----	Hexachlorobenzene	440	U
87-86-5-----	Pentachlorophenol	1100	U
85-01-8-----	Phenanthrene	440	U
120-12-7-----	Anthracene	440	U
86-74-8-----	Carbazole	440	U
84-74-2-----	Di-n-Butylphthalate	1100	U u
206-44-0-----	Fluoranthene	130	J
129-00-0-----	Pyrene	140	J
85-68-7-----	Butylbenzylphthalate	440	U
91-94-1-----	3,3'-Dichlorobenzidine	440	UJ
56-55-3-----	Benzo(a)Anthracene	100	J
218-01-9-----	Chrysene	120	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate	440	U
117-84-0-----	Di-n-Octyl Phthalate	440	U
205-99-2-----	Benzo(b)Fluoranthene	140	J
207-08-9-----	Benzo(k)Fluoranthene	440	U
50-32-8-----	Benzo(a)Pyrene	440	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	440	U
53-70-3-----	Dibenz(a,h)Anthracene	440	U
191-24-2-----	Benzo(g,h,i)Perylene	440	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X207

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344852

Sample wt/vol: 30.60 (g/mL) G Lab File ID: B1102E10

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 26 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/03/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Number TICs found: 20

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q	
1.	UNKNOWN	10.18	1500	BJU	am
2.	UNKNOWN ALIP. ACID	22.37	320	BJU	am
3.	UNKNOWN	23.63	570	J	
4.	UNKNOWN ALIP. HYDROCARBON	23.70	120	J	
5.	UNKNOWN ALIP. HYDROCARBON	24.18	1500	J	
6.	UNKNOWN ALIP. HYDROCARBON	24.92	160	J	
7.	UNKNOWN ALIP. HYDROCARBON	26.13	200	J	
8.	UNKNOWN ALIP. ACID	26.85	150	J	
9.	UNKNOWN ALIP. HYDROCARBON	27.32	160	J	
10.	UNKNOWN ALIP. HYDROCARBON	28.43	130	J	
11.	UNKNOWN ALIP. HYDROCARBON	29.52	200	J	
12.	UNKNOWN ALIP. HYDROCARBON	30.55	310	J	
13.	UNKNOWN ALIP. HYDROCARBON	31.53	440	J	
14.	UNKNOWN ALIP. ACID ESTER	31.57	510	J	
15.	UNKNOWN ALIP. ALCOHOL	32.50	840	J	
16.	UNKNOWN ALIP. HYDROCARBON	33.43	240	J	
17.	UNKNOWN ALIP. HYDROCARBON	34.40	590	J	
18.	UNKNOWN ALIP. HYDROCARBON	36.80	820	J	
19.	UNKNOWN	45.38	790	J	
20.	UNKNOWN	46.15	930	J	

1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X207

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344852

Sample wt/vol: 30.2 (g/mL) G Lab File ID: _____

% Moisture: 26 decanted: (Y/N) N Date Received: 10/27/93

Extraction: (SepF/Cont/Sonc) SONC Date Extracted: 10/28/93

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/04/93

Injection Volume: 1.00 (uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) Y pH: 6.6 Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6	alpha-BHC	2.3	U
319-85-7	beta-BHC	2.3	U
319-86-8	delta-BHC	2.3	U
58-89-9	gamma-BHC (Lindane)	2.3	U
76-44-8	Heptachlor	2.3	U
309-00-2	Aldrin	2.3	0.30 BFPU am
1024-57-3	Heptachlor epoxide	2.3	U
959-98-8	Endosulfan I	2.3	U
60-57-1	Dieldrin	4.4	U
72-55-9	4,4'-DDE	4.4	U
72-20-8	Endrin	4.4	U
33213-65-9	Endosulfan II	4.4	U
50-29-3	4,4'-DDD	4.4	U
1031-07-8	Endosulfan sulfate	4.4	U
50-29-3	4,4'-DDT	4.4	U
72-43-5	Methoxychlor	23	U
53494-70-5	Endrin ketone	4.4	U
7421-36-3	Endrin aldehyde	4.4	U
5103-71-9	alpha-Chlordane	2.3	U
5103-74-2	gamma-Chlordane	2.3	U
8001-35-2	Toxaphene	230	U
12674-11-2	Aroclor-1016	44	U
11104-28-2	Aroclor-1221	90	U
11141-16-5	Aroclor-1232	44	U
53469-21-9	Aroclor-1242	44	U
12672-29-6	Aroclor-1248	44	U
11097-69-1	Aroclor-1254	44	U
11096-82-5	Aroclor-1260	44	U

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X208

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344853

Sample wt/vol: 30.50 (g/mL) G Lab File ID: B1102E06

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 42 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/02/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	560	U
111-44-4	bis(2-Chloroethyl) Ether	560	U
95-57-8	2-Chlorophenol	560	U
541-73-1	1,3-Dichlorobenzene	560	U
106-46-7	1,4-Dichlorobenzene	560	U
95-50-1	1,2-Dichlorobenzene	560	U
95-48-7	2-Methylphenol	560	U
108-60-1	2,2'-oxybis(1-Chloropropane)	560	U
106-44-5	4-Methylphenol	560	U
621-64-7	N-Nitroso-Di-n-Propylamine	560	U
67-72-1	Hexachloroethane	560	U
98-95-3	Nitrobenzene	560	U
78-59-1	Isophorone	560	U
88-75-5	2-Nitrophenol	560	U
105-67-9	2,4-Dimethylphenol	560	U
111-91-1	bis(2-Chloroethoxy)Methane	560	U
120-83-2	2,4-Dichlorophenol	560	U
120-82-1	1,2,4-Trichlorobenzene	560	U
91-20-3	Naphthalene	560	U
106-47-8	4-Chloroaniline	560	UJ
87-68-3	Hexachlorobutadiene	560	U
59-50-7	4-Chloro-3-Methylphenol	560	U
91-57-6	2-Methylnaphthalene	560	U
77-47-4	Hexachlorocyclopentadiene	560	U
88-06-2	2,4,6-Trichlorophenol	560	U
95-95-4	2,4,5-Trichlorophenol	1400	U
91-58-7	2-Chloronaphthalene	560	U
88-74-4	2-Nitroaniline	1400	U
131-11-3	Dimethylphthalate	560	U
208-96-8	Acenaphthylene	560	U
606-20-2	2,6-Dinitrotoluene	560	U
99-09-2	3-Nitroaniline	1400	UJ
83-32-9	Acenaphthene	560	U

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X208

Lab Name: ILLINOIS EPA Contract: 1358070001
 Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846
 Matrix: (soil/water) SOIL Lab Sample ID: D344853
 Sample wt/vol: 30.50 (g/mL) G Lab File ID: B1102E06
 Level: (low/med) LOW Date Received: 10/27/93
 % Moisture: 42 decanted: (Y/N) N Date Extracted: 10/28/93
 Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/02/93
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0
 GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
51-28-5	2,4-Dinitrophenol	1400	U
100-02-7	4-Nitrophenol	1400	U
132-64-9	Dibenzofuran	560	U
121-14-2	2,4-Dinitrotoluene	560	U
84-66-2	Diethylphthalate	560	U
7005-72-3	4-Chlorophenyl-phenylether	560	U
86-73-7	Fluorene	560	U
100-10-6	4-Nitroaniline	1400	U R
534-52-1	4,6-Dinitro-2-methylphenol	1400	U
86-30-6	N-Nitrosodiphenylamine (1)	560	U
101-55-3	4-Bromophenyl-phenylether	560	U
118-74-1	Hexachlorobenzene	560	U
87-86-5	Pentachlorophenol	1400	U
85-01-8	Phenanthrene	560	U
120-12-7	Anthracene	560	U
86-74-8	Carbazole	560	U
84-74-2	Di-n-Butylphthalate	1300	U U
206-44-0	Fluoranthene	560	U
129-00-0	Pyrene	560	U
85-68-7	Butylbenzylphthalate	560	U
91-94-1	3,3'-Dichlorobenzidine	560	UJ
56-55-3	Benzo(a)Anthracene	560	U
218-01-9	Chrysene	560	U
117-81-7	bis(2-Ethylhexyl)Phthalate	560	U
117-84-0	Di-n-Octyl Phthalate	560	U
205-99-2	Benzo(b)Fluoranthene	560	U
207-08-9	Benzo(k)Fluoranthene	560	U
50-32-8	Benzo(a)Pyrene	560	U
193-39-5	Indeno(1,2,3-cd)Pyrene	560	U
53-70-3	Dibenz(a,h)Anthracene	560	U
191-24-2	Benzo(g,h,i)Perylene	560	U

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(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X208

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344853

Sample wt/vol: 30.50 (g/mL) G Lab File ID: B1102E06

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: 42 decanted: (Y/N) N Date Extracted: 10/28/93

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 11/02/93

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 7.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

Number TICs found: 20

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	10.17	2000	BT U
2.	UNKNOWN	10.88	610	BT U
3.	UNKNOWN ALIP. ACID	22.35	460	BT U
4.	UNKNOWN ALIP. HYDROCARBON	24.90	180	J
5.	UNKNOWN ALIP. HYDROCARBON	28.42	140	J
6.	UNKNOWN ALIP. HYDROCARBON	29.50	310	J
7.	UNKNOWN ALIP. HYDROCARBON	30.53	380	J
8.	UNKNOWN ALIP. ACID ESTER	31.57	1100	J
9.	UNKNOWN	32.50	1600	J
10.	UNKNOWN	32.83	170	J
11.	UNKNOWN ALIP. HYDROCARBON	33.42	480	J
12.	UNKNOWN ALIP. HYDROCARBON	34.38	2400	J
13.	UNKNOWN ALIP. HYDROCARBON	35.48	540	J
14.	UNKNOWN	36.03	460	J
15.	UNKNOWN	36.62	610	J
16.	UNKNOWN ALIP. HYDROCARBON	36.80	4800	J
17.	UNKNOWN	36.95	680	J
18.	UNKNOWN ALIP. HYDROCARBON	40.18	1700	J
19.	UNKNOWN	44.75	1500	J
20.	UNKNOWN	47.50	2700	J

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am
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1D
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X208

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344853

Sample wt/vol: 30.8 (g/mL) G Lab File ID: _____

% Moisture: 42 decanted: (Y/N) N Date Received: 10/27/93

Extraction: (SepF/Cont/Sonc) SONC Date Extracted: 10/28/93

Concentrated Extract Volume: 5000 (uL) Date Analyzed: 11/04/93

Injection Volume: 1.00 (uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) Y pH: 7.0 Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6	alpha-BHC	2.9	U
319-85-7	beta-BHC	2.9	U
319-86-8	delta-BHC	2.9	U
58-89-9	gamma-BHC (Lindane)	2.9	U
76-44-8	Heptachlor	2.9	U
309-00-2	Aldrin	2.9	U
1024-57-3	Heptachlor epoxide	2.9	U
959-98-8	Endosulfan I	2.9	U
60-57-1	Dieldrin	1.3	JP
72-55-9	4,4'-DDE	5.5	U
72-20-8	Endrin	2.8	JP
33213-65-9	Endosulfan II	3.6	JP
50-29-3	4,4'-DDD	5.1	J
1031-07-8	Endosulfan sulfate	5.5	U
50-29-3	4,4'-DDT	5.5	U
72-43-5	Methoxychlor	29	U
53494-70-5	Endrin ketone	5.5	U
7421-36-3	Endrin aldehyde	5.5	U
5103-71-9	alpha-Chlordane	0.57	J
5103-74-2	gamma-Chlordane	0.73	JP
8001-35-2	Toxaphene	320	P
12674-11-2	Aroclor-1016	55	U
11104-28-2	Aroclor-1221	110	U
11141-16-5	Aroclor-1232	55	U
53469-21-9	Aroclor-1242	55	U
12672-29-6	Aroclor-1248	55	U
11097-69-1	Aroclor-1254	24	JP
11096-82-5	Aroclor-1260	55	U

2.9 0.59 BJP U am

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

X208RE

Lab Name: ILLINOIS EPA Contract: 1358070001
 Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846
 Matrix: (soil/water) SOIL Lab Sample ID: D344853RE
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK09
 Level: (low/med) LOW Date Received: 10/27/93
 % Moisture: not dec. 42 Date Analyzed: 11/01/93
 GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	Chloromethane	17	U
74-83-9	Bromomethane	17	U
75-01-4	Vinyl Chloride	17	U
75-00-3	Chloroethane	17	U
75-09-2	Methylene Chloride	17	U
67-64-1	Acetone	17	UJ
75-15-0	Carbon Disulfide	17	U
75-35-4	1,1-Dichloroethene	17	U
75-34-3	1,1-Dichloroethane	17	U
540-59-0	1,2-Dichloroethene (total)	17	U
107-06-2	1,2-Dichloroethane	17	U
78-93-3	2-Butanone	17	UJ
71-55-6	1,1,1-Trichloroethane	8	J
56-23-5	Carbon Tetrachloride	17	UJ
75-27-4	Bromodichloromethane	17	UJ
78-87-5	1,2-Dichloropropane	17	UJ
10061-01-5	cis-1,3-Dichloropropene	17	UJ
79-01-6	Trichloroethene	17	UJ
124-48-1	Dibromochloromethane	17	UJ
79-00-5	1,1,2-Trichloroethane	17	UJ
71-43-2	Benzene	17	UJ
10061-02-6	trans-1,3-Dichloropropene	17	UJ
75-25-2	Bromoform	17	UJ
108-10-1	4-Methyl-2-Pentanone	17	UJ
591-78-6	2-Hexanone	17	UJ
127-18-4	Tetrachloroethene	17	UJ
79-34-5	1,1,2,2-Tetrachloroethane	17	UJ
108-88-3	Toluene	17	UJ
108-90-7	Chlorobenzene	17	UJ
100-41-4	Ethylbenzene	17	UJ
100-42-5	Styrene	17	UJ
1330-20-7	Xylene (total)	17	UJ

17 X BU U Am

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

X208RE

Lab Name: ILLINOIS EPA Contract: 1358070001

Lab Code: SPFLD Case No.: EAGLE SAS No.: _____ SDG No.: 344846

Matrix: (soil/water) SOIL Lab Sample ID: D344853RE

Sample wt/vol: 5.0 (g/mL) G Lab File ID: C1101BK09

Level: (low/med) LOW Date Received: 10/27/93

% Moisture: not dec. 42 Date Analyzed: 11/01/93

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

INORGANIC ANALYSIS DATA SHEET

X102

Lab Name: ILLINOIS EPA CHAMPAIGN LAB Contract: EAGLE ZINC CO
 Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: 96
 Matrix (Soil): _____ Lab Sample ID: B316969
 Level (Low/Med): _____ Date Received: 10/28/93
 % Solids: -73.6-

Concentration Units (mg/kg dry weight): _____

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	10000			P
7440-36-0	Antimony	9.2	U	N	P
7440-38-2	Arsenic	5.7		S	FM
7440-39-3	Barium	265			P
7440-41-7	Beryllium	0.81	B		P
7440-43-9	Cadmium	0.61	U		P
7440-70-2	Calcium	9880			P
7440-47-3	Chromium	14.4			P
7440-48-4	Cobalt	6.5	B		P
7440-50-8	Copper	19.7		N	P
7439-89-6	Iron	14400			P
7439-92-1	Lead	236			P
7439-95-4	Magnesium	2090			P
7439-96-5	Manganese	686			P
7439-97-6	Mercury	0.18			AV
7440-02-2	Nickel	11.5			P
7440-09-7	Potassium	1600			P
7782-49-2	Selenium	1.3	U	W	FM
7440-22-4	Silver	1.0	U		P
7440-23-5	Sodium	87.9	B		P
7440-28-0	Thallium	0.34	B	W,N	FM
7440-62-2	Vanadium	27.1			P
7440-66-6	Zinc	138			P
_____	Cyanide	1.1	U		AS
_____	_____	_____	_____	_____	AS
_____	_____	_____	_____	_____	_____

Color Before: -Black- Clarity Before: -Opaque- Texture: -Fine-
 Color After: -Colorless- Clarity After: -Clear- Artifacts: _____
 Comments: _____

INORGANIC ANALYSIS DATA SHEET

X105

Lab Name: ILLINOIS EPA CHAMPAIGN LAB Contract: —EAGLE ZINC CO—
 Lab Code: Case No.: SAS No.: SDG No.: —96—
 Matrix (Soil): Lab Sample ID: —B316960—
 Level (Low/Med): Date Received: 10/28/93
 % Solids: —82.7—

Concentration Units (mg/kg dry weight):

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	7430			P
7440-36-0	Antimony	11.4		N	P
7440-38-2	Arsenic	86.3			FM
7440-39-3	Barium	379			P
7440-41-7	Beryllium	0.83	B		P
7440-43-9	Cadmium	47.2			P
7440-70-2	Calcium	1930			P
7440-47-3	Chromium	22.6			P
7440-48-4	Cobalt	20.1			P
7440-50-8	Copper	911		N	P
7439-89-6	Iron	104000			P
7439-92-1	Lead	5760			P
7439-95-4	Magnesium	1630			P
7439-96-5	Manganese	178			P
7439-97-6	Mercury	0.05	U		AV
7440-02-2	Nickel	55.9			P
7440-09-7	Potassium	300	B		P
7782-49-2	Selenium	1.3		S	FM
7440-22-4	Silver	6.3			P
7440-23-5	Sodium	39.6	B		P
7440-28-0	Thallium	1.3	U	E,N	FM
7440-62-2	Vanadium	22.6			P
7440-66-6	Zinc	31700			P
	Cyanide	1.0	U		AS
					AS

Color Before: —Black— Clarity Before: —Opaque— Texture: —Fine—
 Color After: —Yellow— Clarity After: —Clear— Artifacts:
 Comments:

INORGANIC ANALYSIS DATA SHEET

X107

Lab Name: ILLINOIS EPA CHAMPAIGN LAB Contract: EAGLE ZINC CO
 Lab Code: Case No.: SAS No.: SDG No.: 96
 Matrix (Soil): Lab Sample ID: B316971
 Level (Low/Med): Date Received: 10/28/93
 % Solids: -61.6-

Concentration Units (mg/kg dry weight):

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	13000			P
7440-36-0	Antimony	10.5	U	N	P
7440-38-2	Arsenic	8.7		S	FM
7440-39-3	Barium	124			P
7440-41-7	Beryllium	0.72	B		P
7440-43-9	Cadmium	3.5			P
7440-70-2	Calcium	5360			P
7440-47-3	Chromium	16.1			P
7440-48-4	Cobalt	5.6	Be ₃ V ₃		P
7440-50-8	Copper	36.4		N	P
7439-89-6	Iron	14900			P
7439-92-1	Lead	105			P
7439-95-4	Magnesium	2090			P
7439-96-5	Manganese	600			P
7439-97-6	Mercury	0.16			AV
7440-02-2	Nickel	15.9			P
7440-09-7	Potassium	1160	B		P
7782-49-2	Selenium	1.8	U		FM
7440-22-4	Silver	1.2	U		P
7440-23-5	Sodium	71.8	B		P
7440-28-0	Thallium	0.35	U	W,N	FM
7440-62-2	Vanadium	27.3			P
7440-66-6	Zinc	2480			P
	Cyanide	1.4	U		AS
					AS

Color Before: -Brown Clarity Before: -Opaque Texture: -Fine
 Color After: -Colorless Clarity After: -Clear Artifacts:
 Comments:

INORGANIC ANALYSIS DATA SHEET

X111

Lab Name: ILLINOIS EPA CHAMPAIGN LAB Contract: _____EAGLE ZINC CO_____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: —97—

Matrix (Soil): _____ Lab Sample ID: —B316975—

Level (Low/Med): _____ Date Received: 10/28/93

% Solids: -80.5-

Concentration Units (mg/kg dry weight): _____

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	13500			P
7440-36-0	Antimony	9.0	U	N	P
7440-38-2	Arsenic	8.5			FM
7440-39-3	Barium	193			P
7440-41-7	Beryllium	0.94	B		P
7440-43-9	Cadmium	1.6			P
7440-70-2	Calcium	8380			P
7440-47-3	Chromium	20.2			P
7440-48-4	Cobalt	7.8	B		P
7440-50-8	Copper	33.8			P
7439-89-6	Iron	19600			P
7439-92-1	Lead	70.8			P
7439-95-4	Magnesium	1950			P
7439-96-5	Manganese	491			P
7439-97-6	Mercury	0.11	B		AV
7440-02-2	Nickel	16.5			P
7440-09-7	Potassium	1920			P
7782-49-2	Selenium	0.42	B	W	FM
7440-22-4	Silver	1.0	U		P
7440-23-5	Sodium	120	B		P
7440-28-0	Thallium	0.25	U	W	FM
7440-62-2	Vanadium	34.2			P
7440-66-6	Zinc	488			P
	Cyanide	1.0	U		AS
					AS

Color Before: -Brown- Clarity Before: -Opaque- Texture: -Fine-

Color After: -Colorless- Clarity After: -Clear- Artifacts: _____

Comments: _____

INORGANIC ANALYSIS DATA SHEET

X201

Lab Name: ILLINOIS EPA CHAMPAIGN LAB Contract: _____EAGLE ZINC CO_____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: -97-

Matrix (Soil): _____ Lab Sample ID: -B316978-

Level (Low/Med): _____ Date Received: 10/28/93

% Solids: -69.3-

Concentration Units (mg/kg dry weight): _____

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	6630			P
7440-36-0	Antimony	9.0	U	N	P
7440-38-2	Arsenic	4.5			FM
7440-39-3	Barium	79.5			P
7440-41-7	Beryllium	0.42	B		P
7440-43-9	Cadmium	0.69	B		P
7440-70-2	Calcium	6360			P
7440-47-3	Chromium	9.92			P
7440-48-4	Cobalt	6.1	B		P
7440-50-8	Copper	11.9			P
7439-89-6	Iron	10100			P
7439-92-1	Lead	46.4			P
7439-95-4	Magnesium	2760			P
7439-96-5	Manganese	501			P
7439-97-6	Mercury	0.06	U		AV
7440-02-2	Nickel	9.2	B		P
7440-09-7	Potassium	782	U		P
7782-49-2	Selenium	0.27	U	W	FM
7440-22-4	Silver	0.24			P
7440-23-5	Sodium	73.3	B		P
7440-28-0	Thallium	0.27	U	M	FM
7440-62-2	Vanadium	ca 1790 17.90			P
7440-66-6	Zinc	326			P
	Cyanide	1.2	U		AS
					AS

Color Before: -Brown- Clarity Before: -Opaque- Texture: -Fine-

Color After: -Colorless- Clarity After: -Clear- Artifacts: _____

Comments: _____

INORGANIC ANALYSIS DATA SHEET

X207

Lab Name: ILLINOIS EPA CHAMPAIGN LAB Contract: ~~_____~~EAGLE ZINC CO~~_____~~
 Lab Code: ~~_____~~ Case No.: ~~_____~~ SAS No.: ~~_____~~ SDG No.: ~~_____~~97~~_____~~
 Matrix (Soil): ~~_____~~ Lab Sample ID: ~~_____~~B316984~~_____~~
 Level (Low/Med): ~~_____~~ Date Received: 10/28/93
 % Solids: -74.7-

Concentration Units (mg/kg dry weight): ~~_____~~

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	10700			P
7440-36-0	Antimony	10.7	U	N	P
7440-38-2	Arsenic	6.0			FM
7440-39-3	Barium	167			P
7440-41-7	Beryllium	0.72	B		P
7440-43-9	Cadmium	11.1			P
7440-70-2	Calcium	1510			P
7440-47-3	Chromium	14.6			P
7440-48-4	Cobalt	10.8	B		P
7440-50-8	Copper	20.8			P
7439-89-6	Iron	14900			P
7439-92-1	Lead	76.0			P
7439-95-4	Magnesium	1500			P
7439-96-5	Manganese	1470			P
7439-97-6	Mercury	0.06	U		AV
7440-02-2	Nickel	11.9			P
7440-09-7	Potassium	840	U _B		P
7782-49-2	Selenium	0.33	B	W	FM
7440-22-4	Silver	1.2	U		P
7440-23-5	Sodium	82.0	B		P
7440-28-0	Thallium	0.28	U	W	FM
7440-62-2	Vanadium	41.2			P
7440-66-6	Zinc	2410			P
	Cyanide	1.1	U		AS
					AS

Color Before: ~~_____~~Brown~~_____~~ Clarity Before: ~~_____~~Opaque~~_____~~ Texture: ~~_____~~Fine~~_____~~
 Color After: ~~_____~~Green~~_____~~ Clarity After: ~~_____~~Clear~~_____~~ Artifacts: ~~_____~~
 Comments: ~~_____~~

**IEPA DIVISION OF LABORATORIES
QUALITY ASSURANCE SECTION
INORGANIC DATA VALIDATION
CHECKLIST**

Site: Eagle Zinc Co.
Laboratory: IEPA Champaign
SDG: 97
Analytical Protocol: ILM03.0
Date: December 29, 1993
Reviewer: Chris Bridges
Reviewer Signature: _____

I. PRELIMINARY REVIEW

Number Aqueous Samples: N/A Analytes:
 Number Solid/Soil Samples: 14 Analytes: Trace Metals, Hg, CN

	YES	NO	N/A
A. Chain - of - Custody(ies) Present?	<u>X</u>	_____	_____
Signed ?	<u>X</u>	_____	_____
Dated?	<u>X</u>	_____	_____
B. Cover Page- Present?	<u>X</u>	_____	_____
Do sample numbers agree with sample numbers on:			
a. Chain - of - Custody Forms?	<u>X</u>	_____	_____
b. Form 1s?	<u>X</u>	_____	_____
C. Form 1- Final Data			
Are all Form 1s present and complete?	<u>X</u>	_____	_____
Are correct units indicated on Form 1s (ug/l-waters & mg/kg-soils)	<u>X</u>	_____	_____
Are soil sample results corrected for percent solids (dry weight)?	<u>X</u>	_____	_____
Are sample results < IDL reported as the IDL (U)?	<u>X</u>	_____	_____

ACTIONS:

 NONE

II. HOLDING TIMES & PRESERVATION

Mercury (28 Days)	pH < 2	exceeded?	_____	<u>X</u>	_____
Cyanide (14 Days)	pH > 12	exceeded?	<u>X</u>	_____	_____
other Metals (6months)	pH < 2	exceeded?	_____	<u>X</u>	_____

ACTIONS:

 Cyanide holding time was exceeded. However no data was affected. Aqueous holding times were
 applied to the soil samples

III. CALIBRATIONS

A. Initial Calibration Procedures:

		YES	NO	N/A
Are acceptable 2 point calibrations present for:	ICP?	<u>X</u>	_____	_____
Are acceptable 4 point calibrations present for:	AA?	<u>X</u>	_____	_____
Correlation Coefficient > 0.995?		<u>X</u>	_____	_____
	Cyanide?	<u>X</u>	_____	_____
Correlation Coefficient > 0.995?		<u>X</u>	_____	_____
Mid-Range standard distilled?		<u>X</u>	_____	_____
Are acceptable 4 point calibrations present for:	Mercury?	<u>X</u>	_____	_____
Correlation Coefficient > 0.995?		<u>X</u>	_____	_____
Are acceptable calibrations present for other parameters?		<u>X</u>	_____	_____

ACTIONS:

NONE

B. Form 2 - Initial and Continuing Calibration Verification:

All necessary Form 2s present and complete?	<u>X</u>	_____	_____
ICVs and CCVs analyzed at the correct frequency?	_____	<u>X</u>	_____
Are results reported in the correct units (ug/l)?	<u>X</u>	_____	_____
All calibration verification % Recoveries meet criteria?	<u>X</u>	_____	_____

ACTIONS: (Analyte, % Recovery, Sample(s) affected and Qualifications)

Eleven cyanide samples were analyzed between CCVs however no data was affected

V. ICP INTERFERENCE CHECK SAMPLE:

	YES	NO	N/A
Form 4 present and complete?	<u>X</u>	___	___
Were ICS ran at the correct frequency?	<u>X</u>	___	___
Were all transcription errors corrected?	___	___	<u>X</u>
All % Recoveries of ICSAB Solution +/- 20 % of True Value?	<u>X</u>	___	___
For elements not present in ICESA, is the absolute value of the ICESA result greater than the IDL?	<u>X</u>	___	___

ACTIONS: (Analyte, % Recovery, Sample(s) affected, Qualifications)

The results for Co, Ni, V, and Mn were all greater than their respective IDLs however no data was affected

VI. SPIKE SAMPLE RECOVERY:

Form 5 present and complete for:	each 20 samples?	<u>X</u>	___	___
	each matrix type?	___	___	<u>X</u>
Were all transcription errors corrected?		___	___	<u>X</u>
Were field blanks used for spike sample analysis?		___	<u>X</u>	___
Were all Matrix Spike % Recoveries within criteria?		___	<u>X</u>	___

ACTIONS: (Analyte, % Recovery, Sample(s) affected, Qualifications)

Sb (46.9%R) X206 is qualified as estimated (J); X108, X109, X110, X111, X112, X113, X201, X202, X203, X204, X205, X206, X207, X208 is qualified as estimated (UJ)

Pb by GFAA (175.1 %R) the sample result was > 4X the spike added so no recovery limits apply

IX. FURNACE ATOMIC ABSORPTION (AA) QC:

	YES	NO	N/A
Did the laboratory utilize duplicate injections for all non-MSA analyses?	<u>X</u>	___	___
Does the GFAA flow chart appear to have been followed for all analyses?	<u>X</u>	___	___
Did the laboratory properly flag the GFAA results on the Form 1s?	<u>X</u>	___	___

ACTIONS: (Analyte, Sample(s) affected, Qualifications)

Results flagged 'W', 'E', or 'M' by the laboratory are qualified as estimated (J)

X. ICP SERIAL DILUTION:

Form 9 present and complete?	<u>X</u>	___	___
Was Serial Dilution analysis performed for:			
each 20 or fewer samples	<u>X</u>	___	___
each matrix type?	___	___	<u>X</u>
Were all transcription errors corrected?	___	___	<u>X</u>
Were all serial dilution results within criteria?	<u>X</u>	___	___
Were field blanks used for serial dilution analysis?	___	<u>X</u>	___

ACTIONS: (Analyte, Sample(s) affected, Qualifications)

NONE

XI. RAW DATA:

	YES	NO	N/A
Digestion Log for flame AA/ICP present?	<u>X</u>	_____	_____
Digestion Log for furnace AA present?	_____	_____	<u>X</u>
Digestion Log for mercury present?	<u>X</u>	_____	_____
Digestion Log for cyanide present?	<u>X</u>	_____	_____
Inventory Sheet present?	<u>X</u>	_____	_____
Weights, dilutions, and volumes used to obtain values present?	<u>X</u>	_____	_____
Percent solids calculation present for soils (sediments)?	<u>X</u>	_____	_____
Are preparation dates present on Digestion Logs?	<u>X</u>	_____	_____
Are standards preparation logs present and dated?	<u>X</u>	_____	_____
Measurement read out records present for:			
ICP?	<u>X</u>	_____	_____
Flame AA?	_____	_____	<u>X</u>
Furnace AA?	<u>X</u>	_____	_____
Mercury?	<u>X</u>	_____	_____
Cyanide?	<u>X</u>	_____	_____
other Inorganics?	_____	_____	<u>X</u>
Are all results with the ICP linear ranges?	<u>X</u>	_____	_____
Are all Mercury results within the calibrated range?	<u>X</u>	_____	_____
Are all Cyanide results within the calibrated range?	<u>X</u>	_____	_____
Are all other inorganics within the calibrated range?	_____	_____	<u>X</u>
Sulfate, Sulfide, _____			
Are all raw data to support all sample analyses an QC operations present?	<u>X</u>	_____	_____
Legible?	<u>X</u>	_____	_____
Properly labeled?	<u>X</u>	_____	_____
Are the Instrument Detection Limits (IDLs) less than 3 months old?	<u>X</u>	_____	_____
Are the ICP Linear Ranges less than three months old?	<u>X</u>	_____	_____
Are the ICP Interelement correction factors less than one year old?	<u>X</u>	_____	_____

ACTIONS:

NONE

**IEPA DIVISION OF LABORATORIES
QUALITY ASSURANCE SECTION
INORGANIC DATA VALIDATION
CHECKLIST**

Site: Eagle Zinc (1 of 2)
Laboratory: IEPA Champaign
SDG: 96
Analytical Protocol: ILMO3.0
Date: December 21, 1993
Reviewer: Chris Bridges
Reviewer Signature: *Chris Bridges*

I. PRELIMINARY REVIEW

Number Aqueous Samples: N/A Analytes:
 Number Solid/Soil Samples: 14 Analytes: Trace Metals, Hg, CN

	YES	NO	N/A
A. Chain - of - Custody(ies) Present?	<u>X</u>	_____	_____
Signed ?	<u>X</u>	_____	_____
Dated?	<u>X</u>	_____	_____
B. Cover Page- Present?	<u>X</u>	_____	_____
Do sample numbers agree with sample numbers on:			
a. Chain - of - Custody Forms?	<u>X</u>	_____	_____
b. Form 1s?	<u>X</u>	_____	_____
C. Form 1- Final Data			
Are all Form 1s present and complete?	<u>X</u>	_____	_____
Are correct units indicated on Form 1s (ug/l-waters & mg/kg-soils)	<u>X</u>	_____	_____
Are soil sample results corrected for percent solids (dry weight)?	<u>X</u>	_____	_____
Are sample results < IDL reported as the IDL (U)?	<u>X</u>	_____	_____
D. Data Package received within the required turn around time?	_____	<u>X</u>	_____

ACTIONS:

NONE

II. HOLDING TIMES & PRESERVATION

Mercury (28 Days)	pH < 2	exceeded?	_____	<u>X</u>	_____
Cyanide (14 Days)	pH > 12	exceeded?	<u>X</u>	_____	_____
other Metals (6months)	pH < 2	exceeded?	_____	<u>X</u>	_____
Contactual Holding Times?		exceeded?	<u>X</u>	_____	_____

ACTIONS:

Cyanides technical holding time exceeded by 4 days however, no data is affected.

III. CALIBRATIONS

A. Initial Calibration Procedures:

	YES	NO	N/A
Are acceptable 2 point calibrations present for:			
ICP?	<u>X</u>	___	___
Are acceptable 4 point calibrations present for:			
AA?	<u>X</u>	___	___
Correlation Coefficient > 0.995?	<u>X</u>	___	___
Cyanide?	<u>X</u>	___	___
Correlation Coefficient > 0.995?	<u>X</u>	___	___
Mid-Range standard distilled?	<u>X</u>	___	___
Are acceptable 4 point calibrations present for:			
Mercury?	<u>X</u>	___	___
Correlation Coefficient > 0.995?	<u>X</u>	___	___
Are acceptable calibrations present for other parameters?	<u>X</u>	___	___

ACTIONS:

NONE

B. Form 2 - Initial and Continuing Calibration Verification:

All necessary Form 2s present and complete?	<u>X</u>	___	___
ICVs and CCVs analyzed at the correct frequency?	<u>X</u>	___	___
Are results reported in the correct units (ug/l)?	<u>X</u>	___	___
All calibration verification % Recoveries meet criteria?	<u>X</u>	___	___

ACTIONS: (Analyte, % Recovery, Sample(s) affected and Qualifications)

NONE

V. ICP INTERFERENCE CHECK SAMPLE:

	YES	NO	N/A
Form 4 present and complete?	<u>X</u>	___	___
Were ICS run at the correct frequency?	<u>X</u>	___	___
Were all transcription errors corrected?	___	___	<u>X</u>
All % Recoveries of ICSAB Solution +/- 20 % of True Value?	<u>X</u>	___	___
For elements not present in ICESA, is the absolute value of the ICESA result greater than the IDL?	<u>X</u>	___	___

ACTIONS: (Analyte, % Recovery, Sample(s) affected, Qualifications)

The results for Cr, Cu, Mn, and Zn all exceeded their respective IDLs however no data was affected

VI. SPIKE SAMPLE RECOVERY:

Form 5 present and complete for:	each 20 samples?	<u>X</u>	___	___
	each matrix type?	<u>X</u>	___	___
Were all transcription errors corrected?		___	___	<u>X</u>
Were field blanks used for spike sample analysis?		___	<u>X</u>	___
Were all Matrix Spike % Recoveries within criteria?		___	<u>X</u>	___

ACTIONS: (Analyte, % Recovery, Sample(s) affected, Qualifications)

Sb (31.4 %R) X105 is qualified as estimated (J); X103, X104, X114, X115, X116, X117, X118, X119
X120, X101, X102, X106, X107 are qualified as estimated (UJ)

Cu (126.9 %R) X103, X104, X105, X114, X115, X116, X117, X118, X119, X120, X101, X102, X106,
X107 are qualified as estimated (J)

Tl (71.6 %R) X103, X104, X105, X116, X117, X118, X119, X120, X107 are qualified as estimated (UJ)
X101, X102, X106, X114, X115 are qualified as estimated (J)

VII. DUPLICATE SAMPLE ANALYSIS:

	YES	NO	N/A
Form 6 present and complete for: each 20 samples?	<u>X</u>	___	___
each matrix type?	___	___	<u>X</u>
Were all transcription errors corrected?	___	___	<u>X</u>
Were field blanks used for duplicate analysis?	___	<u>X</u>	___
For both AA and ICP when both are used for the same analyte?	<u>X</u>	___	___
Were all duplicate analyses differences within criteria?	<u>X</u>	___	___

ACTIONS: (Element, Differences, Sample(s) affected, Qualifications)

NONE

VIII. LABORATORY CONTROL SAMPLE:

(Note: LCS not required for aqueous Hg.)

Form 7 Present and Complete?	<u>X</u>	___	___
Was one LCS prepared and analyzed for:			
every 20 or fewer water samples?	___	___	<u>X</u>
every digestion batch of water samples?	___	___	<u>X</u>
every 20 or fewer solid samples?	<u>X</u>	___	___
every digestion batch of solid samples?	<u>X</u>	___	___
Were all transcription errors corrected?	___	___	<u>X</u>
Were all of the Aqueous LCS % Recoveries within criteria?	___	___	<u>X</u>
Were all of the Solid LCS % Recoveries within criteria?	<u>X</u>	___	___

ACTIONS: (Element, % Recovery, Sample(s) affected, Qualifications)

NONE

IX. FURNACE ATOMIC ABSORPTION (AA) QC:

	YES	NO	N/A
Did the laboratory utilize duplicate injections for all non-MSA analyses?	<u>X</u>	___	___
Does the GFAA flow chart appear to have been followed for all analyses?	<u>X</u>	___	___
Did the laboratory properly flag the GFAA results on the Form 1s?	<u>X</u>	___	___

ACTIONS: (Analyte, Sample(s) affected, Qualifications)

Any result flagged 'W' or 'E' by the laboratory is qualified as estimated (J)

X. ICP SERIAL DILUTION:

Form 9 present and complete?	<u>X</u>	___	___
Was Serial Dilution analysis performed for:			
each 20 or fewer samples	<u>X</u>	___	___
each matrix type?	___	___	<u>X</u>
Were all transcription errors corrected?	___	___	<u>X</u>
Were all serial dilution results within criteria?	<u>X</u>	___	___
Were field blanks used for serial dilution analysis?	___	<u>X</u>	___

ACTIONS: (Analyte, Sample(s) affected, Qualifications)

NONE

XI. RAW DATA:

	YES	NO	N/A
Digestion Log for flame AA/ICP present?	<u>X</u>	_____	_____
Digestion Log for furnace AA present?	<u>X</u>	_____	_____
Digestion Log for mercury present?	<u>X</u>	_____	_____
Digestion Log for cyanide present?	<u>X</u>	_____	_____
Inventory Sheet present?	<u>X</u>	_____	_____
Weights, dilutions, and volumes used to obtain values present?	<u>X</u>	_____	_____
Percent solids calculation present for soils (sediments)?	<u>X</u>	_____	_____
Are preparation dates present on Digestion Logs?	<u>X</u>	_____	_____
Are standards preparation logs present and dated?	<u>X</u>	_____	_____
Measurement read out records present for:			
ICP?	<u>X</u>	_____	_____
Flame AA?	_____	_____	<u>X</u>
Furnace AA?	<u>X</u>	_____	_____
Mercury?	<u>X</u>	_____	_____
Cyanide?	<u>X</u>	_____	_____
other Inorganics?	_____	_____	<u>X</u>
Are all results with the ICP linear ranges?	<u>X</u>	_____	_____
Are all Mercury results within the calibrated range?	<u>X</u>	_____	_____
Are all Cyanide results within the calibrated range?	<u>X</u>	_____	_____
Are all other inorganics within the calibrated range?	_____	_____	<u>X</u>
Sulfate, Sulfide, _____			
Are all raw data to support all sample analyses an QC operations present?	<u>X</u>	_____	_____
Legible?	<u>X</u>	_____	_____
Properly labeled?	<u>X</u>	_____	_____
Are the Instrument Detection Limits (IDLs) less than 3 months old?	<u>X</u>	_____	_____
Are the ICP Linear Ranges less than three months old?	<u>X</u>	_____	_____
Are the ICP Interelement correction factors less than one year old?	<u>X</u>	_____	_____

ACTIONS: NONE

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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PRELIMINARY REVIEW

1. Chain-of-Custody

YES NO

- a. [] Check chain-of-custody documentation for date/time sampled, date/time received in laboratory.
- b. [] Check chain-of-custody documentation for proper documentation of transfers and signoffs.
- c. [] Check chain-of-custody documentation for any inconsistencies or anomalies.

Comments:

None.

2. Case Narrative

YES NO

- a. [] Review entire case narrative.
- b. [] Check case narrative for completeness.
- c. [] Check for proper authorization signature.

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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I. Holding Times

YES NO

[] Check that all technical and/or contractual holding times were met, as required, for all fractions.

EPA Number	Lab Number	Date Coll.	Date Rec'd.	VOA	Semi-VOA		Pesticide	
				Date Anal.	Date Extr	Date Anal.	Date Extr	Date Anal.
X201	-846	10/26/93	10/27/93	10/29/93	10/28/93	11/3/93	10/28/93	11/4/93
X202	-847	↓	↓	↓	↓	11/3	↓	↓
X203	-848	↓	↓	↓	↓	11/3	↓	↓
X204	-849	↓	↓	↓	↓	11/3	↓	↓
X205	-850	↓	↓	↓	↓	11/3	↓	↓
X206	-851	↓	↓	↓	↓	11/2	↓	↓
X207	-852	↓	↓	↓	↓	11/3	↓	↓
X208	-853	↓	↓	↓	↓	11/2	↓	↓
X203RE				11/01/93				
X204RE				↓				
X205RE								
X206RE								
X207RE				↓				
X208RE								

List below all samples (by sample number and fraction) qualified due to holding times.

All holding times met.

Data Validation Checklist

Site Name: Eagle Fire Co.

SDG

No.: 344846

Laboratory: IEPA

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II. GC/MS Instrument Performance Check

Fraction: VOA SemiVOA (circle one)

1. Evaluate Forms V and Raw Data

- | | YES | NO | |
|----|-------------------------------------|--------------------------|------------------------------------------------------------------------------------------------------------------------|
| a. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check that Forms V are present and completed for each 12 hour time period. |
| b. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for transcription errors between raw data and Forms V. |
| c. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check that the appropriate number of significant figures has been reported and that rounding errors have not occurred. |
| d. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for calculation errors. |

2. Verify Raw Data Format

- | | YES | NO | |
|--|-------------------------------------|--------------------------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check mass spectral listing to determine that the mass assignment is correct and that the mass listing is normalize to the specified ion (m/z 95 for VOA, m/z 198 for SemiVOA). |

3. Verify Ion Abundance Criteria

- | | YES | NO | |
|--|-------------------------------------|--------------------------|-----------------------------------------------------|
| | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check that all ion abundance criteria has been met. |

4. Verify Background Correction

- | | YES | NO | |
|--|-------------------------------------|--------------------------|--------------------------------------------------------------------------------------------|
| | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check that tuning compound spectra were generated using appropriate background correction. |

Comments: None.

VOLATILE CALIBRATION OUTLIERS

3/90 SOW

Lab Name: IEPA

Case: Eagle Zinc Co.

Instrument #	Minimum RRF	Initial Cal.			Contin. Cal.			Contin. Cal.			Contin. Cal.		
		RF	%RSD	Q									
XL2		10/28/93	1117		10/29/93	933		11/01/93	1038				
heated purge													
Chloromethane	0.010		37.8										
Bromomethane	0.100												
Vinyl Chloride	0.100												
Chloroethane	0.010												
Methylene Chloride	0.010												
Acetone	0.010							32.9	J				
Carbon Disulfide	0.010												
1,1-Dichloroethene	0.100												
1,1-Dichloroethane	0.200												
1,2-Dichloroethene (total)	0.010												
Chloroform	0.200												
1,2-Dichloroethane	0.100												
2-Butanone	0.010				29.4	J		49.4	J				
1,1,1-Trichloroethane	0.100												
Carbon Tetrachloride	0.100												
Bromodichloromethane	0.200												
1,2-Dichloropropane	0.010												
cis-1,3-Dichloropropene	0.200												
1,1,2-Dichloroethane	0.300												
Dibromochloromethane	0.100												
1,1,2-Trichloroethane	0.100												
Benzene	0.500												
trans-1,3-Dichloropropene	0.100												
Bromoform	0.100												
4-Methyl-2-Pentanone	0.010							25.4	J				
2-Hexanone	0.010				25.4	J		35.0	J				
Tetrachloroethene	0.200												
1,1,2,2-Tetrachloroethane	0.500												
Toluene	0.400												
Chlorobenzene	0.500												
Ethylbenzene	0.100												
Styrene	0.300												
Xylene(total)	0.300												
Bromofluorobenzene	0.300												

AFFECTED SAMPLES:

Reviewers Initials/Date
am
 12/9/93

	VBLKSL1	VBLKSL2
	X201	X203RF
	X202	X204RF
	X203	X205RF
	X204	X206RF
	X205	X207RF
	X206	X208RF
	X207	X203MS
	X208	X203MSD

This column of flags should be applied to the analytes on the sample data sheets.

Data Validation Checklist

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: IEPA

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III. Initial Calibration

GC/MS

Fraction: VOA SemiVOA (circle one)

1. Verify that the Correct Standard Concentrations Were Used.

YES NO

[] Check the Forms VI and the raw data to verify that the correct standard concentrations were used to calibrate the GC/MS instrument(s).

2. Verify that the Correct Initial Calibration was Used for Water and Low Level Soils.

YES NO N/A

[] [] Check that initial calibrations were performed as required for water/med. level soil and low level soil.

3. Verify Use of Correct Standards.

YES NO N/A

[] [] Check that the correct standard was used for quantitation of samples, if samples were analyzed immediately subsequent to initial calibration.

4. Evaluate Initial Calibration RRFs and \overline{RRF} s.

YES NO

a. [] Check and recalculate the RRFs and \overline{RRF} s for several target compounds (at least one associated with each internal standard).

b. [] Check that, for all target compounds and surrogates, the \overline{RRF} s meet the applicable criteria. Note any "outliers" on the Calibration Outliers Form.

5. Evaluate Initial Calibration %RSDs.

YES NO

a. [] Check and recalculate the %RSD for several target compounds.

b. [] Check that the applicable %RSD criteria have been met. Note any "outliers" on the Calibration Outliers Form.

Comments: See Calibration Outliers Form.

Data Validation Checklist

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: EPA

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IV. Continuing Calibration

GC/MS

Fraction: (VOA) SemiVOA (circle one)

1. Verify Continuing Calibration Frequency.

YES NO

[] Check the continuing calibration raw data and Forms VII to verify that continuing calibration standards were analyzed at the proper frequency and that each continuing calibration was compared to the appropriate initial calibration.

2. Evaluate Continuing Calibration RRFs.

YES NO

a. [] Check and recalculate the continuing calibration RRFs for several compounds.

b. [] Check that all target compound and surrogate RRFs meet the criteria.

3. Evaluate Continuing Calibration %Ds.

YES NO

a. [] Check and recalculate the continuing calibration %Ds for several compounds.

b. [] Check that all target compound and surrogate %Ds meet the applicable criteria.

Comments:

See Calibration Outlier's Form.

Data Validation Checklist

Site Name: Eagle Zinn Co

SDG

No.: 344846

Laboratory: IEPA

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VI. Surrogate Spikes

GC/MS

Fraction: VOA SemiVOA (circle one)

1. Review Raw Data.

YES NO

[] Check raw data to verify that the recoveries on the Form II are accurate and within the limits.

2. Evaluate Surrogate Recovery Calculations.

YES NO

[] Check that the surrogate spike recoveries were calculated correctly and are free from transcription errors.

3. Evaluate Surrogate Recoveries.

YES NO

a. [] Check that reanalyses were performed as required.

b. [] Check that surrogate recoveries in blanks met criteria.

4. Evaluate Reanalyses.

YES NO

[] Whenever there are two or more analyses for a particular sample, determine which are the best analyses to use. This determination must be performed in conjunction with the evaluation of the internal standard area response criteria. List below the results of the reviewers determinations.

Comments:

X205 + X207 - BFB high ? R
X206 + X206RE - Toluene-d8 high ? R

4. Use X203RE, X204RE, X205RE, X206RE, X207, X208RE

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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VII. Matrix Spikes/Matrix Spike Duplicates

Fraction: VOA SemiVOA Pesticide (circle one)

1. Verify Frequency

YES NO

[] Check that MS and MSD samples were analyzed at the correct frequency.

2. Evaluate MS/MSD Criteria.

YES NO

[] Check MS/MSD results for %R and RPD are within the advisory limits.

3. Verify MS/MSD Calculations.

YES NO

a. [] Check raw data and verify that results are calculated correctly and are free from transcription errors.

b. [] Check that %Rs and RPDs were calculated correctly.

4. Evaluate Sample Precision.

YES NO

[] Compare %RSD results of non-spiked compounds between the original result, MS and MSD.

X 203 Compound	Orig. Result	MS Result	MSD Result	%RSD
MeCl ₂	9	6	6	25%
Acetone	15	10	8	33%

Comments:

None

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: LEPA

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X. Internal Standards

GC/MS

Fraction: VOA SemiVOA (circle one)

1. Evaluate Raw Data.

YES NO

[] Check raw data and verify that the internal standard retention times and areas reported on the Forms VIII are correct.

2. Verify RT and IS Area Criteria.

YES NO

[] Check that retention times and internal standard area meet the appropriate criteria.

3. Evaluate Reanalyses.

YES NO

[] Whenever there are two or more analyses for a particular sample, determine which are the best analyses to use. This determination must be performed in conjunction with the evaluation of the surrogate spike recovery criteria. List the results of the reviewers determinations in Section VI., Surrogate Spikes.

Comments:

X203, X204, X204RE, X205, X205RE, X206, X206RE, ~~X207~~ X208,
+ X208RE - DFB + CB2 low area

X203RE, X207RE, X203MS, X203MSD - DFB low area

Data Validation Checklist

Site Name: Eagle zinc Co

SDG

No.: 344 P46

Laboratory: IEPA

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XI. Target Compound Identification

GC/MS

Fraction: (VOA) SemiVOA (circle one)

1. Verify Relative Retention Time (RRT) Criteria.

YES NO

[] Check that the RRT of reported compounds is within the criteria.

2. Evaluate Target Compound Spectra.

YES NO

[] Check the sample target compound spectra against the laboratory standard spectra; verify that the specified criteria are met.

3. Evaluate Possible Carryover.

YES NO

[] Check the raw data of the samples as related to the samples analyzed previously to verify that sample carryover has not adversely affected results.

4. Evaluate Chromatograms.

YES NO

[] Check the sample chromatograms to verify that peaks are accounted for.

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: LEPA

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XII. Compound Quantitation and Reported CRQLs

Fraction: VOA SemiVOA Pesticide (circle one)

1. Evaluate Quantitation of Sample Results.

YES NO

[] Check raw data to verify calculation of sample results.

2. Evaluate Quantitation Parameters.

YES NO N/A

[] []

For GC/MS analyses, check that the correct internal standard, quantitation ion, and *RRF* were used to quantitate results. Verify that the same internal standard, quantitation ion, and *RRF* are used throughout, in both the calibration and as well as the quantitation process.

3. Evaluate CRQLs.

YES NO

[] Check that the CRQLs have been adjusted to reflect all sample dilutions, concentrations, splits, cleanup activities, and dry weight factors.

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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XIII. Tentatively Identified Compounds

GC/MS Only

Fraction: VOA SemiVOA (circle one)

1. Evaluate Tentative Identifications.

YES NO

[] Check that all TICs reported meet the identification guidelines.

2. Evaluate Raw Data.

YES NO

[] Check raw data to verify that the laboratory has generated a library search for all required peaks in the chromatograms for samples and blanks.

3. Evaluate Blanks.

YES NO

[] Check blank sample chromatograms to verify that TIC peaks present in samples are not found in blanks.

4. Examine Mass Spectra.

YES NO

[] Check all mass spectra for every sample.

5. Evaluate TIC Identifications.

YES NO

[] Since TIC library searches often yield several candidate compounds, all reasonable choices must be considered.

6. Evaluate Laboratory Artifacts and Contaminants.

YES NO

[] Check sample results and raw data to verify that common laboratory artifacts and contaminants are not reported as sample contaminants.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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XIII. TICs continued

7. Evaluate Possibility of False Negatives.

YES NO N/A
a. [] []

Check to determine if target compounds have been identified and quantitated as TICs.

b. [] []

If target compounds have been identified and quantitated as TICs, check to determine whether the false negative is an isolated occurrence or whether additional data may be affected. Comment on all such false negatives below.

8. Determine That Results Are From Proper Fraction.

YES NO N/A
 [] []

Target compounds could be identified in more than one fraction; if this occurs, check that quantitation is from the proper fraction.

9. Verify That Internal Standards And Surrogates Are Not Searched.

YES NO
 []

Check that library searches were not performed on internal standards or surrogates.

10. Verify Estimated Quantitation of TICs.

YES NO
 []

Check that the estimated concentration of the TICs was made using an assumed RRF of one.

Comments:

None

Data Validation Checklist

Site Name: Eagle Iron Co.

SDG

No.: 344846

Laboratory: - LEPA

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XIV. GC/MS System Performance

Fraction: VOA SemiVOA (circle one)

1. Evaluate Overall System Performance.

- | | YES | NO | |
|----|-------------------------------------|--------------------------|---------------------------------------------------------------------------------------------------|
| a. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for high RIC background levels or shifts in absolute retention times of internal standards. |
| b. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for excessive baseline rise at elevated temperature. |
| c. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for extraneous peaks. |
| d. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for loss of resolution. |
| e. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for peak tailing or peak splitting that may result in inaccurate quantitation. |

Comments:

None

Data Validation Checklist

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: EPA

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II. GC/MS Instrument Performance Check

Fraction: VOA SemiVOA (circle one)

1. Evaluate Forms V and Raw Data

- YES NO
- a. [] Check that Forms V are present and completed for each 12 hour time period.
- b. [] Check for transcription errors between raw data and Forms V.
- c. [] Check that the appropriate number of significant figures has been reported and that rounding errors have not occurred.
- d. [] Check for calculation errors.

2. Verify Raw Data Format

- YES NO
- [] Check mass spectral listing to determine that the mass assignment is correct and that the mass listing is normalize to the specified ion (m/z 95 for VOA, m/z 198 for semiVOA).

3. Verify Ion Abundance Criteria

- YES NO
- [] Check that all ion abundance criteria has been met.

4. Verify Background Correction

- YES NO
- [] Check that tuning compound spectra were generated using appropriate background correction.

Comments: None.

Data Validation Checklist

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: IEPA

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III. Initial Calibration

GC/MS

Fraction: VOA SemiVOA (circle one) (circle one)

1. Verify that the Correct Standard Concentrations Were Used.

YES NO

[] Check the Forms VI and the raw data to verify that the correct standard concentrations were used to calibrate the GC/MS instrument(s).

2. Verify that the Correct Initial Calibration was Used for Water and Low Level Soils.

YES NO N/A

[] [] Check that initial calibrations were performed as required for water/med. level soil and low level soil.

3. Verify Use of Correct Standards.

YES NO N/A

[] [] Check that the correct standard was used for quantitation of samples, if samples were analyzed immediately subsequent to initial calibration.

4. Evaluate Initial Calibration RRFs and \overline{RRF} s.

YES NO

a. [] Check and recalculate the RRFs and \overline{RRF} s for several target compounds (at least one associated with each internal standard).

b. [] Check that, for all target compounds and surrogates, the \overline{RRF} s meet the applicable criteria. Note any "outliers" on the Calibration Outliers Form.

5. Evaluate Initial Calibration %RSDs.

YES NO

a. [] Check and recalculate the %RSD for several target compounds.

b. [] Check that the applicable %RSD criteria have been met. Note any "outliers" on the Calibration Outliers Form.

Comments:

See Calibration Outliers Form.

Data Validation Checklist

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: IEPA

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IV. Continuing Calibration

GC/MS

Fraction: VOA (SemiVOA) (circle one)

1. Verify Continuing Calibration Frequency.

YES NO

[] Check the continuing calibration raw data and Forms VII to verify that continuing calibration standards were analyzed at the proper frequency and that each continuing calibration was compared to the appropriate initial calibration.

2. Evaluate Continuing Calibration RRFs.

YES NO

a. [] Check and recalculate the continuing calibration RRFs for several compounds.

b. [] Check that all target compound and surrogate RRFs meet the criteria.

3. Evaluate Continuing Calibration %Ds.

YES NO

a. [] Check and recalculate the continuing calibration %Ds for several compounds.

b. [] Check that all target compound and surrogate %Ds meet the applicable criteria.

Comments:

See Calibration Duties Form.

SEMIVOLATILE CALIBRATION OUTLIERS

Page 1

Lab Name: LEPA

Case: Eagle zinc Co.

Instrument # / NC 500	Minimum RRF	Initial Cal.			Contin. Cal.			Contin. Cal.			Contin. Cal.		
		DATE/TIME:	RF	%RSD	Q	RF	%RSD	Q	RF	%RSD	Q	RF	%RSD
		9/14/93	1242		11/2/93	1939		11/3/93	1255				
Phenol	0.800												
bis(2-Chloroethyl)ether	0.700												
2-Chlorophenol	0.800												
1,3-Dichlorobenzene	0.600												
1,4-Dichlorobenzene	0.500												
1,2-Dichlorobenzene	0.400												
2-Methylphenol	0.700												
2,2'-oxybis(1-Chloropropanol)	0.010												
4-Methylphenol	0.600												
N-Nitroso-di-n-propylamine	0.500												
Hexachloroethane	0.300												
Nitrobenzene	0.200												
Isoclorane	0.400												
2-Nitrophenol	0.100												
2,4-Dimethylphenol	0.200												
bis(2-Chloroethoxy)methane	0.300												
2,4-Dichlorophenol	0.200												
1,2,4-Trichlorobenzene	0.200												
Phthalene	0.700												
4-Chloroaniline	0.010	70.1			60.2	J							
Hexachlorobutadiene	0.010												
4-Chloro-3-methylphenol	0.200												
2-Methylnaphthalene	0.400												
Hexachlorocyclopentadiene	0.010												
2,4,5-Trichlorophenol	0.200												
2,4,5-Trichlorophenol	0.200												
2-Chloronaphthalene	0.800												
2-Nitroaniline	0.010												
Dimethylphthalate	0.010												
Acenaphthylene	1.300												
2,5-Dinitrotoluene	0.200												
3-Nitroaniline	0.010	46.9			70.9	J							
Acenaphthene	0.800												
2,4-Dinitrophenol	0.010												
4-Nitrophenol	0.010												
Dibenzofuran	0.800												
2,4-Dinitrotoluene	0.200												

AFFECTED SAMPLES:

Reviewer's Initials/Date

am
12/10/93

	SBLKSL	X208MS
	X206	X208MSD
	X208	
	X205	
	X202	
	X201	
	X207	
	X203	
	X204	

SEMIVOLATILE CALIBRATION OUTLIERS

Page 2

Name: LEPACase: Eagle Zinc

Instrument #	Minimum	Initial Cal.			Contn. Cal.			Contn. Cal.			Contn. Cal.		
		DATE	TIME	RRF	RF	%RSD	Q	RF	%RSD	Q	RF	%RSD	Q
INC500		9/14/93	1242		11/2/93	1935		11/3/93	1255				
Diethylthalate	0.010												
4-Chlorophenyl-phenylether	0.400												
Fluorene	0.900												
4-Nitroaniline	0.010		90.6		.015	74.6	J	.026	55.9	J			
4,6-Dinitro-2-methylphenol	0.010												
N-Nitrosodiphenylamine (1)	0.010												
4-Bromophenyl-ethylether	0.100												
Hexachlorobenzene	0.100												
Pentachlorophenol	0.050												
Phenanthrene	0.700												
Anthracene	0.700												
Carbazole	0.010												
Di-n-butylthalate	0.010												
Fluoranthene	0.500												
Pyrene	0.600												
Butylbenzylthalate	0.010												
3,3'-Dichlorobenzidine	0.010					31.2	J						
Benzo(a)anthracene	0.800												
Chrysene	0.700												
bis(2-Ethylhexyl)phthalate	0.010												
-octylthalate	0.010												
Benzo(b)fluoranthene	0.700												
Benzo(k)fluoranthene	0.700												
Benzo(a)pyrene	0.700												
Indeno(1,2,3-cd)pyrene	0.500												
Dibenz(a,h)anthracene	0.400												
Benzo(g,h,i)perylene	0.500												
Nitrobenzene-d5	0.200												
2-Fluorobiphenyl	0.700												
Terphenyl-d14	0.500												
Phenol-d6	0.800												
2-Fluorophenol	0.600												
2-Chlorophenol-d4	0.800												
1,2-Dichlorobenzene-d4	0.400												
4-Nitrophenol	0.010												
Dibenzofuran	0.800												
2,4-Dinitrotoluene	0.200												

3/92

Q - This column of flags should be applied to the analytes on the sample data sheets.

SEE PAGE 1 FOR AFFECTED SAMPLES

Reviewer's
Initials/Date

am
12/10/93

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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VII. Matrix Spikes/Matrix Spike Duplicates

Fraction: VOA SemiVOA Pesticide (circle one)

1. Verify Frequency

YES NO

[] Check that MS and MSD samples were analyzed at the correct frequency.

2. Evaluate MS/MSD Criteria.

YES NO

[] Check MS/MSD results for %R and RPD are within the advisory limits.

3. Verify MS/MSD Calculations.

YES NO

a. [] Check raw data and verify that results are calculated correctly and are free from transcription errors.

b. [] Check that %Rs and RPDs were calculated correctly.

4. Evaluate Sample Precision.

YES NO

[] Compare %RSD results of non-spiked compounds between the original result, MS and MSD.

X208	Compound	Orig. Result	MS Result	MSD Result	%RSD
	Di-n-Butyl phthalate	1300	1400	1500	7%

Comments:

X208MS + X208MSD - 4-Nitrophenol + Pentachlorophenol high %R

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: LEPA

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X. Internal Standards

GC/MS

Fraction: VOA SemiVOA (circle one) (circle one)

1. Evaluate Raw Data.

YES NO

[] Check raw data and verify that the internal standard retention times and areas reported on the Forms VIII are correct.

2. Verify RT and IS Area Criteria.

YES NO

[] Check that retention times and internal standard area meet the appropriate criteria.

3. Evaluate Reanalyses.

YES NO

[] Whenever there are two or more analyses for a particular sample, determine which are the best analyses to use. This determination must be performed in conjunction with the evaluation of the surrogate spike recovery criteria. List the results of the reviewers determinations in Section VI., Surrogate Spikes.

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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XI. Target Compound Identification

GC/MS

Fraction: VOA SemiVOA (circle one)

1. Verify Relative Retention Time (RRT) Criteria.

YES NO

[] Check that the RRT of reported compounds is within the criteria.

2. Evaluate Target Compound Spectra.

YES NO

[] Check the sample target compound spectra against the laboratory standard spectra; verify that the specified criteria are met.

3. Evaluate Possible Carryover.

YES NO

[] Check the raw data of the samples as related to the samples analyzed previously to verify that sample carryover has not adversely affected results.

4. Evaluate Chromatograms.

YES NO

[] Check the sample chromatograms to verify that peaks are accounted for.

Comments:

None.

Data Validation Checklist.

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: TEPA

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XII. Compound Quantitation and Reported CRQLs

Fraction: VOA SemiVOA Pesticide (circle one)

1. Evaluate Quantitation of Sample Results.

YES NO

[] Check raw data to verify calculation of sample results.

2. Evaluate Quantitation Parameters.

YES NO N/A

[] []

For GC/MS analyses, check that the correct internal standard, quantitation ion, and *RRF* were used to quantitate results. Verify that the same internal standard, quantitation ion, and *RRF* are used throughout, in both the calibration and as well as the quantitation process.

3. Evaluate CRQLs.

YES NO

[] Check that the CRQLs have been adjusted to reflect all sample dilutions, concentrations, splits, cleanup activities, and dry weight factors.

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinic Co

SDG

No.: 344846

Laboratory: IEPA

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XIII. Tentatively Identified Compounds

GC/MS Only

Fraction: VOA SemiVOA (circle one)

1. Evaluate Tentative Identifications.

YES NO

[] Check that all TICs reported meet the identification guidelines.

2. Evaluate Raw Data.

YES NO

[] Check raw data to verify that the laboratory has generated a library search for all required peaks in the chromatograms for samples and blanks.

3. Evaluate Blanks.

YES NO

[] Check blank sample chromatograms to verify that TIC peaks present in samples are not found in blanks.

4. Examine Mass Spectra.

YES NO

[] Check all mass spectra for every sample.

5. Evaluate TIC Identifications.

YES NO

[] Since TIC library searches often yield several candidate compounds, all reasonable choices must be considered.

6. Evaluate Laboratory Artifacts and Contaminants.

YES NO

[] Check sample results and raw data to verify that common laboratory artifacts and contaminants are not reported as sample contaminants.

Data Validation Checklist

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: LEDA

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XIII. TICs continued

7. Evaluate Possibility of False Negatives.

YES NO N/A
a. [] []

Check to determine if target compounds have been identified and quantitated as TICs.

b. [] []

If target compounds have been identified and quantitated as TICs, check to determine whether the false negative is an isolated occurrence or whether additional data may be affected. Comment on all such false negatives below.

8. Determine That Results Are From Proper Fraction.

YES NO N/A
 [] []

Target compounds could be identified in more than one fraction; if this occurs, check that quantitation is from the proper fraction.

9. Verify That Internal Standards And Surrogates Are Not Searched.

YES NO
 []

Check that library searches were not performed on internal standards or surrogates.

10. Verify Estimated Quantitation of TICs.

YES NO
 []

Check that the estimated concentration of the TICs was made using an assumed RRF of one.

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co
SDG

No.: 344846

Laboratory: - IEDPA

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XIV. GC/MS System Performance

Fraction: VOA SemiVOA (circle one)

1. Evaluate Overall System Performance.

- | | YES | NO | |
|----|-------------------------------------|--------------------------|---------------------------------------------------------------------------------------------------|
| a. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for high RIC background levels or shifts in absolute retention times of internal standards. |
| b. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for excessive baseline rise at elevated temperature. |
| c. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for extraneous peaks. |
| d. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for loss of resolution. |
| e. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check for peak tailing or peak splitting that may result in inaccurate quantitation. |

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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II. Pesticide Instrument Performance Check

1. Resolution Check Mixture

- | | YES | NO | |
|----|-------------------------------------|--------------------------|------------------------------------------------------------------------------------------------------------------|
| a. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the Form VIII PEST. to determine that the resolution check mixture(s) was analyzed in the proper sequence. |
| b. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the resolution check mixture data and the Form VI PEST.-4 to verify that the resolution criterion was met. |

2. Performance Evaluation Mixture

- | | YES | NO | |
|----|-------------------------------------|--------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------|
| a. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the Form VII PEST. to determine that the PEM(s) was analyzed at the proper frequency and position in the initial calibration sequence. |
| b. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the PEM data from the initial and continuing calibrations to verify that the resolution criterion was met. |
| c. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the PEM data from the initial and continuing calibrations and Form VII PEST.-1 to verify that the retention times are within the retention time windows. |
| d. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check that the <i>RPDs</i> meet the criterion. |
| e. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check that the breakdowns for 4,4'-DDT and Endrin meet the criteria. |

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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III. Initial Calibration

PESTICIDES

1. Individual Standard Mixtures.

- | | YES | NO | |
|----|-------------------------------------|--------------------------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| a. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the Form VIII PEST to verify that the Individual Standard Mixtures were analyzed at the proper frequency for each GC column and instrument. Check that the proper concentrations were used. |
| b. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the raw data to determine that the midpoint standard is at the proper concentration and verify that the resolution criterion has been met for each midpoint concentration standard. |
| c. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the Individual Standard Mixture data and Form VI PEST.-1 and review the calculated retention time windows for calculation and transcription errors. |
| d. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the Individual Standard Mixture data and Form VI PEST.-2 to verify that the %RSDs for the calibration factors meet the criterion. Check and recalculate several %RSDs for errors. |

2. Multi Component Compounds.

- | | | | |
|----|-------------------------------------|--------------------------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| a. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the raw data and the Form VIII PEST. to verify that the Multi-component Standards were analyzed at the proper concentration and frequency for each GC column and instrument. |
| b. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the raw data and Form VI PEST.-3 to verify that at least three peaks were used for calibration and that retention time and calibration factor data are available for each peak. |

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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IV. Continuing Calibration

PESTICIDES

1. Evaluate Continuing Calibration Standards.

YES NO

[] Check the Form VIII PEST to verify that the Instrument Blanks, PEMs, and Individual Standard Mixtures were analyzed at the proper frequency and that no more than 12:00 hours elapsed between calibration brackets in an ongoing analytical sequence.

2. Individual Standard Mixtures Resolution.

YES NO

[] Check the data for the midpoint concentration of the Individual Standard Mixtures to verify that the resolution criteria was met.

3. Individual Standard Mixtures Retention Times

YES NO

[] Check the data for each of the single component pesticides and surrogates in the midpoint concentration of the Individual Standard Mixtures to verify that the retention times are within the appropriate windows.

4. Evaluate Continuing Calibration RPDs.

YES NO

[] Check the data for the midpoint concentration of the Individual Standard Mixtures and Form VII PEST.-2 to verify that the RPDs between the calculated amount and the true amount for each of the pesticides and surrogates meet the criterion.

Comments:

4. IndAMO3 (col DB-1701) 10/29/93 1648
Methoxychlor 26.5% RPD - high * No samples affected

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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V. Blanks

Fraction: VOA SemiVOA

Pest. (circle one)

1. Review Blank Results.

YES NO

[] Check all associated blanks for the presence of TCL compounds or TICs. Note all contaminated blanks and associated samples below.

2. Verify Blank Frequency.

YES NO

[] Check that blank analyses have been performed at the required frequency.

Blank Summary

Blank Sample No. PBLK51 PBLK12 _____
 Date Anal. or Extr. 10/28/93 Ext. 11/5/93 Analyzed _____
 Instrument HP 1+2 HP 1+2 _____

TCL Comp'd.	Amount						
Aldrin	0.37	DDT	0.0052				
		Toxaphene	0.12				

TIC Comp'd.	Amount						
NA		NA					

Comments: All other instrument blanks and GPC blank are O.K.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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VI. Surrogate Spikes

Pesticides

1. Review Raw Data.

YES NO

[] Check raw data to verify that the recoveries on the Form II are accurate and within the limits.

2. Evaluate Surrogate Recovery Calculations.

YES NO

[] Check that the surrogate spike recoveries were calculated correctly and are free from transcription errors.

3. Evaluate Possible Interferences.

YES NO N/A

[] [] If surrogate spike recoveries are not acceptable, check the raw data for possible interferences which may have effected surrogate recoveries.

4. Evaluate Retention Times.

YES NO N/A

[] [] If retention time limits are not met, check the raw data for possible misidentification of GC peaks.

5. Evaluate Any Low Recoveries.

YES NO N/A

[] [] If low surrogate recoveries are observed, check whether low recoveries are due to sample dilution.

6. Evaluate Surrogate Analyses in Blanks.

YES NO

[] Check that all surrogate analysis criteria (retention time and advisory recovery criteria) were met in all blank samples.

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: LEPA

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VII. Matrix Spikes/Matrix Spike Duplicates

Fraction: VOA SemiVOA Pesticide (circle one)

1. **Verify Frequency**

YES NO
 [] Check that MS and MSD samples were analyzed at the correct frequency.

2. **Evaluate MS/MSD Criteria.**

YES NO
 [] Check MS/MSD results for %R and RPD are within the advisory limits.

3. **Verify MS/MSD Calculations.**

- a. YES NO
 [] Check raw data and verify that results are calculated correctly and are free from transcription errors.
- b. [] Check that %Rs and RPDs were calculated correctly.

4. **Evaluate Sample Precision.**

YES NO
 [] Compare %RSD results of non-spiked compounds between the original result, MS and MSD.

X205	Compound	Orig. Result	MS Result	MSD Result	%RSD
	Endrin ketone	1.6	3.2	2.5	33%

Comments:

None

Data Validation Checklist

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: IEPA

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X. Pesticide Cleanup Checks

1. Florisil Cartridge Check.

- YES NO
- a. [] Check the data from the Florisil cartridge solution analyses and the Form IX PEST.-1 and check some of the %R calculations; verify that there are no calculation or transcription errors.
- b. [] Check all criteria have been met.

2. Gel Permeation Chromatography.

- YES NO
- a. [] Check the data from the GPC calibration check analyses and the Form IX PEST.-2 and recalculate some of the %R results; verify that there are no calculation or transcription errors.
- b. [] Check all criteria have been met and that Arochlor patterns are similar to those of previous standards.

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co

SDG

No.: 344846

Laboratory: IEPA

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XI. Target Compound Identification

Pesticides

1. Evaluate Reported Results.

- | | YES | NO | |
|----|-------------------------------------|--------------------------|-------------------------------------------------------------------------------------------------------------------------------------|
| a. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the Form I PEST., the associated raw data, and Form X PEST.-1 and Form X PEST.-2 to confirm reported detected analytes. |
| b. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the Form I PEST., the associated raw data, and Form X PEST.-1 and Form X PEST.-2 to confirm reported non-detects. |
| c. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the associated blank data for potential interferences to evaluate sample data for false positives. |
| d. | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Check the calibration data for adequate retention time windows to evaluate the sample data for false positives and false negatives. |

2. Evaluate Multi-Component Analyte Results.

- | | YES | NO | |
|--|-------------------------------------|--------------------------|------------------------------------------------------------------------------------------------------------------------------------------------|
| | <input checked="" type="checkbox"/> | <input type="checkbox"/> | Compare the retention times and relative peak height ratios of major multi-component analyte peaks against appropriate standard chromatograms. |

3. Verify GC/MS Confirmations if Applicable.

- | | YES | NO | N/A | |
|--|--------------------------|--------------------------|-------------------------------------|-------------------------------------------------------------------------------------------------------------------------------|
| | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | Check that GC/MS confirmation was performed for pesticide concentrations in the final sample extract which exceeded 10 ng/ul. |

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: IEPA

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XII. Compound Quantitation and Reported CRQLs

Fraction: VOA SemiVOA Pesticide (circle one)

1. Evaluate Quantitation of Sample Results.

YES NO

[] Check raw data to verify calculation of sample results.

2. Evaluate Quantitation Parameters.

YES NO N/A

[] []

For GC/MS analyses, check that the correct internal standard, quantitation ion, and RRF were used to quantitate results. Verify that the same internal standard, quantitation ion, and RRF are used throughout, in both the calibration and as well as the quantitation process.

3. Evaluate CRQLs.

YES NO

[] Check that the CRQLs have been adjusted to reflect all sample dilutions, concentrations, splits, cleanup activities, and dry weight factors.

Comments:

None.

Data Validation Checklist

Site Name: Eagle Zinc Co.

SDG

No.: 344846

Laboratory: LEPA

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XV. Overall Assessment of Data

Evaluate the Overall Quality of the Data.

YES NO

[] Evaluate any technical problems which have not been previously addressed.

[] Review all available materials to assess the overall quality of the data, keeping in mind the additive nature of analytical problems.

[] If appropriate information is available, assess the usability of the data to assist the data user in avoiding inappropriate use of the data. Review all available information, including the QAPjP, SAP, and communications with the data user that concerns the intended use of the data.

Provide a brief narrative to give the data user an indication of the analytical limitations of the data. Include any details from the above checks. Any inconsistency of the data with the Case Narrative should be noted. If sufficient information is available, the reviewer should give an assessment of the usability of the data within the given context.

Data are useable as qualified on Forms 1.

Alicia Mudd

Multiple horizontal lines for additional text or signature.